LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Title : SVOA CALIBRATION

Signal : TIC: BG051351.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.655	364	369	375	rVV	193366	313386	1.66%	0.343%
2	5.790	385	392	403	rBV	521602	1101820	5.83%	1.205%
3	6.161	448	455	464	rVB2	272653	607567	3.21%	0.664%
4	6.648	525	538	543	rBV3	63560	161946	0.86%	0.177%
5	7.142	618	622	627	rVV2	178273	297988	1.58%	0.326%
6	7.195	627	631	636	rVV	117170	177414	0.94%	0.194%
7	7.253	636	641	646	rVV	230742	389817	2.06%	0.426%
8	7.312	646	651	657	rVV3	229288	409855	2.17%	0.448%
9	7.389	657	664	671	rVB2	368249	720513	3.81%	0.788%
10	7.559	687	693	698	rBV	95392	182869	0.97%	0.200%
11	7.776	725	730	734	rVV	622750	949007	5.02%	1.038%
12	7.823	734	738	746	rVB	450235	754702	3.99%	0.825%
13	8.135	786	791	796	rVB	182463	276956	1.47%	0.303%
14	8.194	796	801	804	rBV	98519	193936	1.03%	0.212%
15	8.229	804	807	813	rVB	135007	228790	1.21%	0.250%
16	8.299	813	819	824	rBV2	149021	280704	1.49%	0.307%
17	8.687	880	885	889	rVV3	117122	257822	1.36%	0.282%
18	8.746	889	895	902	rVB2	158067	400854	2.12%	0.438%
19	8.822	902	908	920	rBV2	330555	620115	3.28%	0.678%
20	8.928	920	926	930	rBV	149741	242403	1.28%	0.265%
21	9.286	976	987	997	rBV3	153205	354034	1.87%	0.387%
22	9.398	997	1006	1011	rBV	861622	1464711	7.75%	1.602%
23	9.539	1026	1030	1037	rVV4	81701	193133	1.02%	0.211%
24	9.633	1040	1046	1059	rVV7	84556	267399	1.41%	0.292%
25	9.815	1072	1077	1083	rVV2	105877	207085	1.10%	0.226%
26	9.880	1083	1088	1092	rVV2	95949	176503	0.93%	0.193%
27	10.185	1135	1140	1144	rVV4	77126	162385	0.86%	0.178%
28	10.397	1170	1176	1182	rVV3	160045	291927	1.54%	0.319%
29	10.949	1264	1270	1277	rVV	526230	884348	4.68%	0.967%
30	11.026	1277	1283	1285	rVV	151697	292642	1.55%	0.320%
31	11.078	1285	1292	1297	rVV	1160324	2255428	11.93%	2.466%
32	11.131	1297	1301	1307	rVV	135177	230143	1.22%	0.252%
33	11.995	1440	1448	1452	rBV	144585	256978	1.36%	0.281%
34	12.389	1509	1515	1522	rBV	518644	777637	4.11%	0.850%
35	12.665	1556	1562	1569	rVB	652546	1094317	5.79%	1.197%
36	12.882	1592	1599	1607	rVB	303746	560197	2.96%	0.613%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Title : SVOA CALIBRATION

37	13.276	1660	1666	1671	rBV	188612	316992	1.68%	0.347%
38	13.329	1671	1675	1680	rVB	133444	187573	0.99%	0.205%
39	13.617	1717	1724	1728	rBV	756813	1120418	5.93%	1.225%
40	13.858	1760	1765	1775	rVB	84290	172177	0.91%	0.188%
41	14.004	1782	1790	1796	rVB3	157113	407212	2.15%	0.445%
42	14.145	1808	1814	1817	rVV2	233231	484553	2.56%	0.530%
43	14.175	1817	1819	1827	rVV2	213522	529639	2.80%	0.579%
44	14.263	1831	1834	1839	rVV2	192397	309665	1.64%	0.339%
45	14.386	1850	1855	1858	rVV	124148	194712	1.03%	0.213%
46	14.439	1860	1864	1873	rVB4	81316	186876	0.99%	0.204%
47	14.680	1901	1905	1911	rVB	715483	904495	4.79%	0.989%
48	14.827	1926	1930	1935	rVV	221677	340591	1.80%	0.372%
49	15.021	1958	1963	1971	rBV2	89723	231501	1.22%	0.253%
50	15.221	1992	1997	2004	rVV4	225490	422803	2.24%	0.462%
51	15.291	2004	2009	2017	rVB3	201350	343844	1.82%	0.376%
52	15.432	2028	2033	2038	rBV2	94738	199574	1.06%	0.218%
53	15.591	2054	2060	2061	rBV3	151068	251034	1.33%	0.275%
54	15.632	2062	2067	2072	rVB	678613	972361	5.14%	1.063%
55	15.738	2081	2085	2090	rBV2	123330	189841	1.00%	0.208%
56	16.031	2131	2135	2139	rVV	214006	347715	1.84%	0.380%
57	16.078	2139	2143	2148	rVB3	159764	275879	1.46%	0.302%
58	16.496	2209	2214	2224	rBV	662874	1550932	8.21%	1.696%
59	16.625	2231	2236	2240	rBV2	209542	348821	1.85%	0.381%
60	16.772	2255	2261	2264	rBV5	111487	220439	1.17%	0.241%
61	17.283	2344	2348	2352	rBV	557683	698547	3.70%	0.764%
62	17.336	2353	2357	2361	rVB	224493	322329	1.71%	0.352%
63	17.442	2371	2375	2379	rVB	158903	218998	1.16%	0.239%
64	17.583	2394	2399	2403	rBV	259854	396346	2.10%	0.433%
65	17.829	2437	2441	2446	rVB	165814	241066	1.28%	0.264%
66	18.023	2470	2474	2481	rVB	505038	695918	3.68%	0.761%
67	18.452	2543	2547	2553	rBV	434176	676355	3.58%	0.740%
68	18.511	2553	2557	2563	rVB	290021	435109	2.30%	0.476%
69	18.711	2588	2591	2595	rVB	366306	436117	2.31%	0.477%
70	18.993	2636	2639	2647	rVB3	191522	288292	1.53%	0.315%
71	19.339	2693	2698	2703	rBV2	557239	792230	4.19%	0.866%
72	19.539	2728	2732	2737	rBV2	256674	354883	1.88%	0.388%
73	19.768	2768	2771	2776	rVB	260537	306709	1.62%	0.335%
74	19.821	2777	2780	2784	rVB	159296	190939	1.01%	0.209%
75	19.909	2791	2795	2798	rBV	303725	372884	1.97%	0.408%
76	19.939	2798	2800	2807	rVB2	300497	512652	2.71%	0.561%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Title : SVOA CALIBRATION

77	20.221	2844	2848	2851	rBV	295211	377721	2.00%	0.413%
78	20.438	2881	2885	2890	rBV	558457	681023	3.60%	0.745%
79	20.879	2951	2960	2964	rVB	9244492	17912762	94.77%	19.589%
80	20.937	2966	2970	2974	rVB	510623	591229	3.13%	0.647%
81	21.196	3011	3014	3020	rVB2	489030	703489	3.72%	0.769%
82	21.360	3038	3042	3049	rBV	172891	274960	1.45%	0.301%
83	21.460	3055	3059	3065	rVB	923547	1336592	7.07%	1.462%
84	21.742	3098	3107	3120	rVB	9301215	18902007	100.00%	20.670%
85	21.860	3122	3127	3129	rBV3	237292	431232	2.28%	0.472%
86	22.024	3149	3155	3160	rBV2	637886	1132065	5.99%	1.238%
87	22.483	3230	3233	3235	rBV	151269	205952	1.09%	0.225%
88	22.530	3235	3241	3247	rVB4	1022994	2157010	11.41%	2.359%
89	22.659	3258	3263	3274	rVB2	600553	1138362	6.02%	1.245%
90	22.994	3314	3320	3335	rVB	686525	1776248	9.40%	1.942%
91	23.158	3343	3348	3356	rBV4	193018	437064	2.31%	0.478%
92	23.382	3380	3386	3397	rVB2	446972	958119	5.07%	1.048%
93	23.881	3466	3471	3481	rVB	218422	472290	2.50%	0.516%
94	24.228	3524	3530	3538	rBV	300600	651197	3.45%	0.712%
95	24.633	3593	3599	3612	rVB	436992	1608508	8.51%	1.759%
96	24.909	3640	3646	3664	rVB4	352405	1039498	5.50%	1.137%
97	25.221	3693	3699	3705	rBV2	209354	452308	2.39%	0.495%
98	26.413	3895	3902	3913	rVB2	180802	560618	2.97%	0.613%
99	26.913	3979	3987	4002	rVB2	282697	958311	5.07%	1.048%
100	27.113	4012	4021	4032	rVB5	344777	1200353	6.35%	1.313%

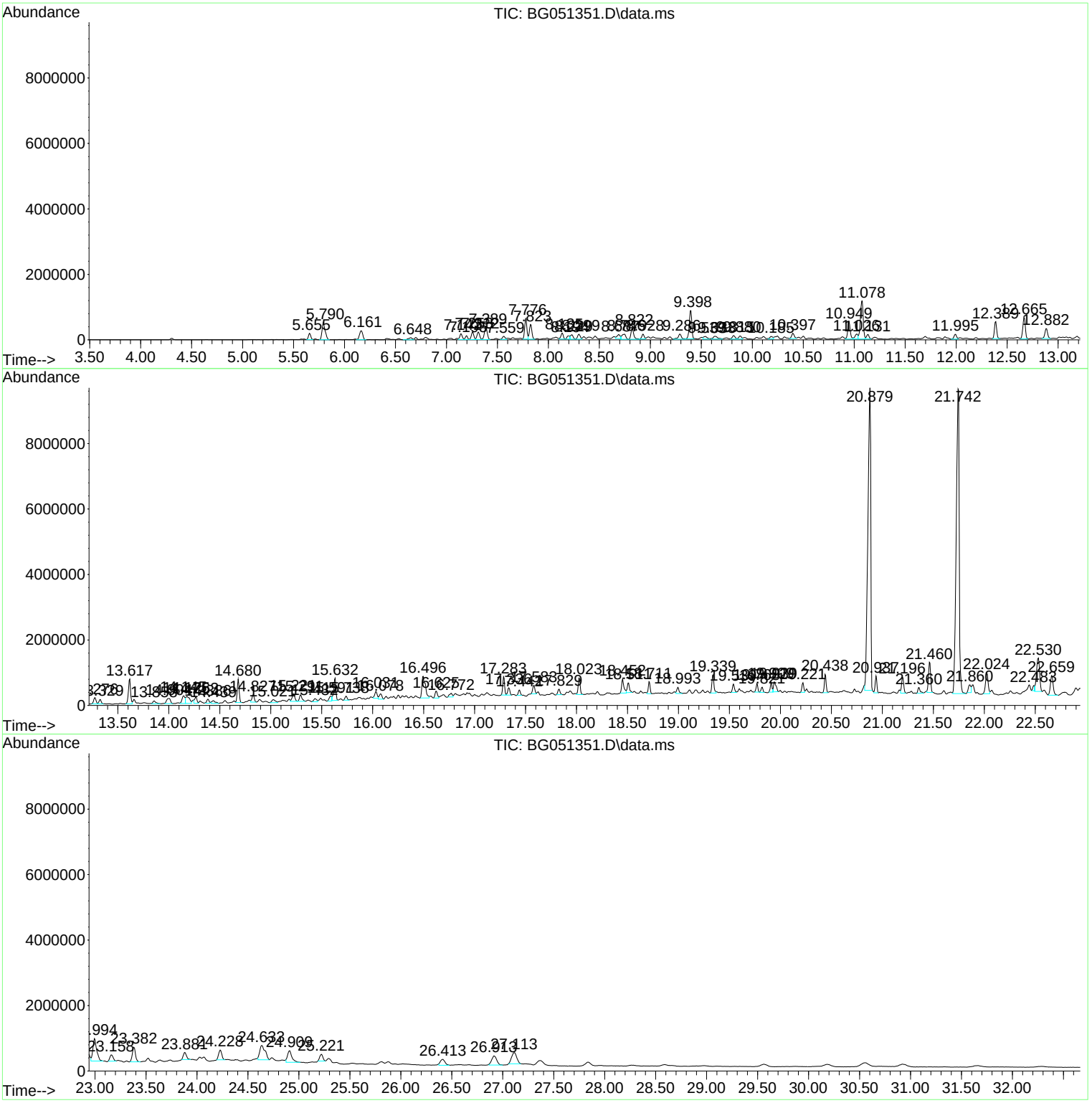
Sum of corrected areas: 91445270

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

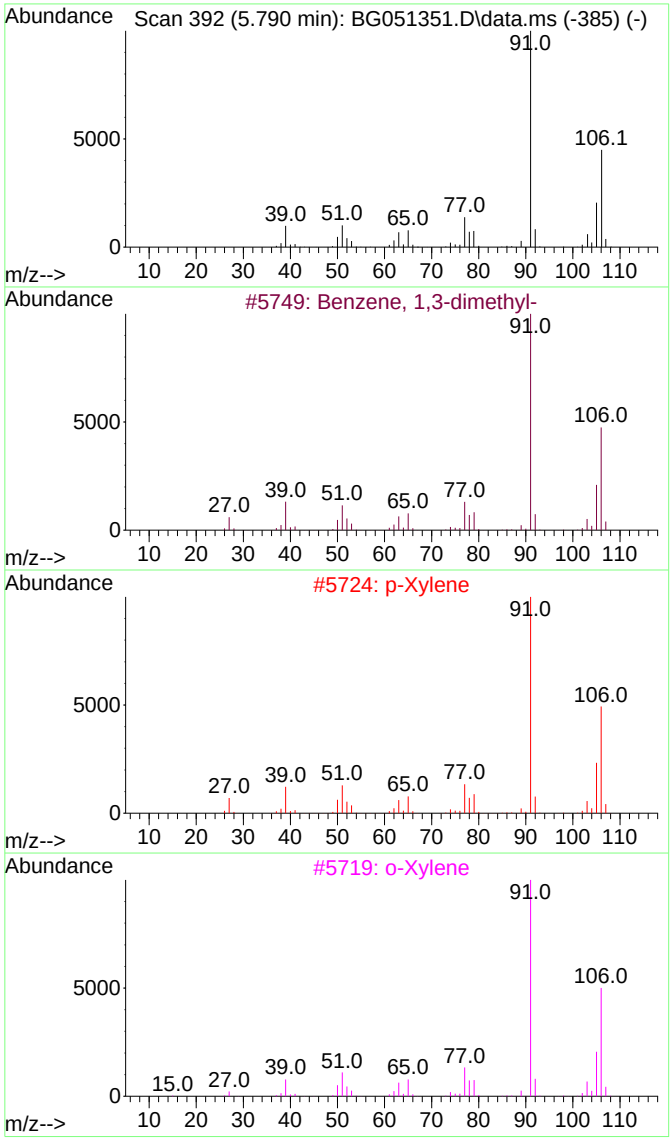
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

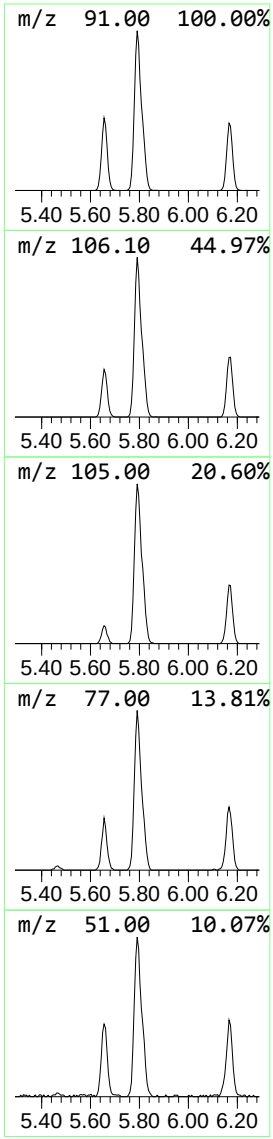
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.790	113.63 ng/ul	1101820	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2	p-Xylene	106	C8H10	000106-42-3	97
3	o-Xylene	106	C8H10	000095-47-6	97
4	p-Xylene	106	C8H10	000106-42-3	97
5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

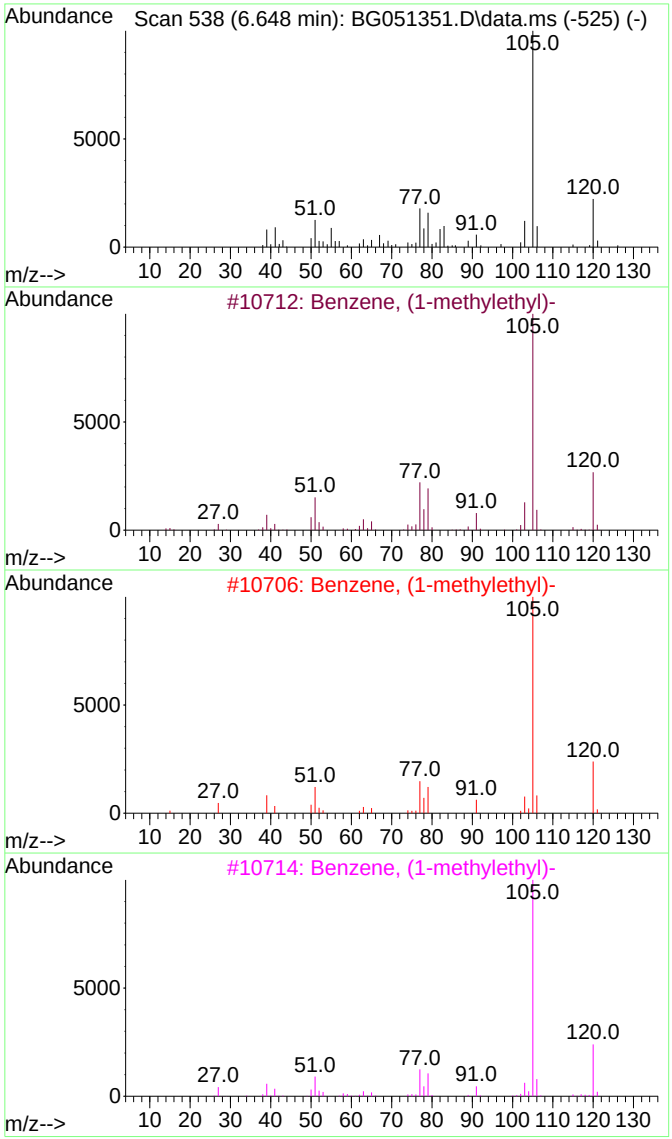
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

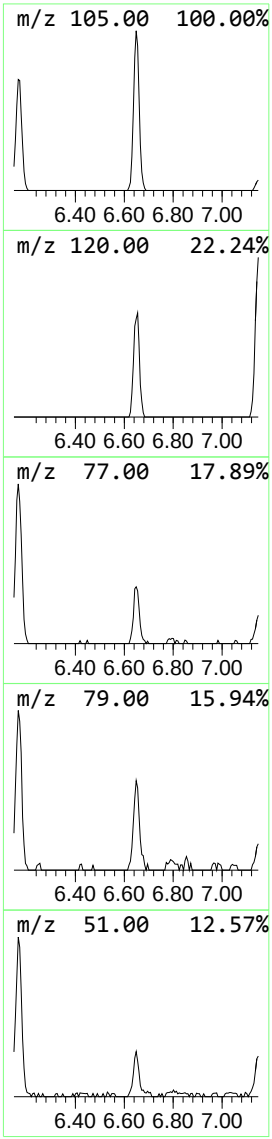
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, (1-methylethyl)- Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.648	16.70 ng/ul	161946	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	95
2	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	93
3	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	90
4	Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	87
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

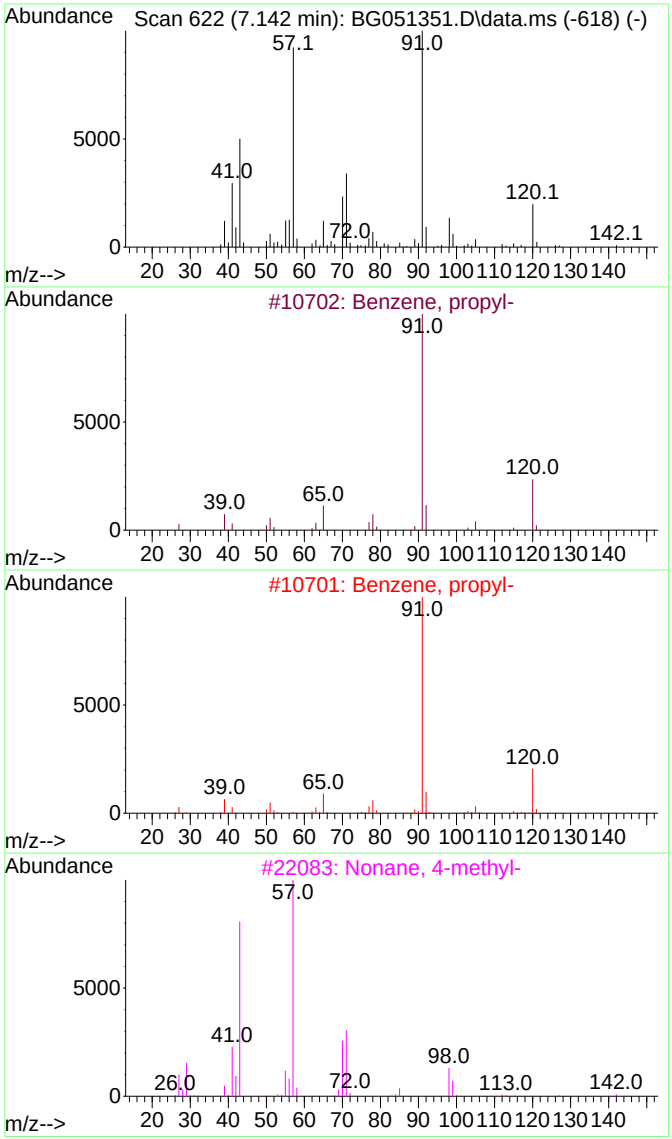
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

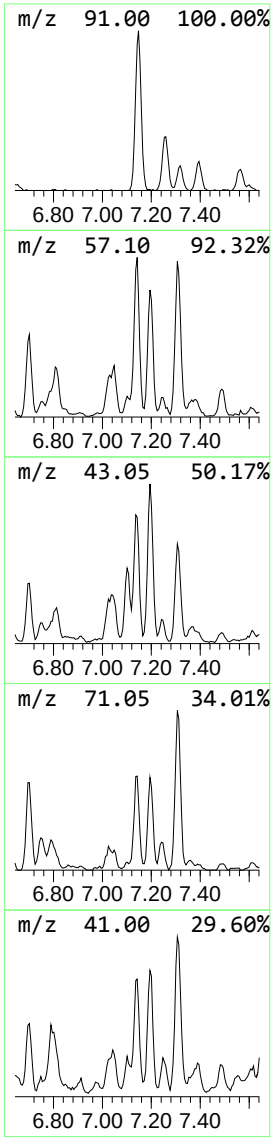
****

Peak Number 5 unknown-01 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.142	30.73 ng/ul	297988	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	46
2	Benzene, propyl-	120	C9H12	000103-65-1	46
3	Nonane, 4-methyl-	142	C10H22	017301-94-9	45
4	Benzeneacetaldehyde	120	C8H8O	000122-78-1	42
5	Benzene, propyl-	120	C9H12	000103-65-1	41





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

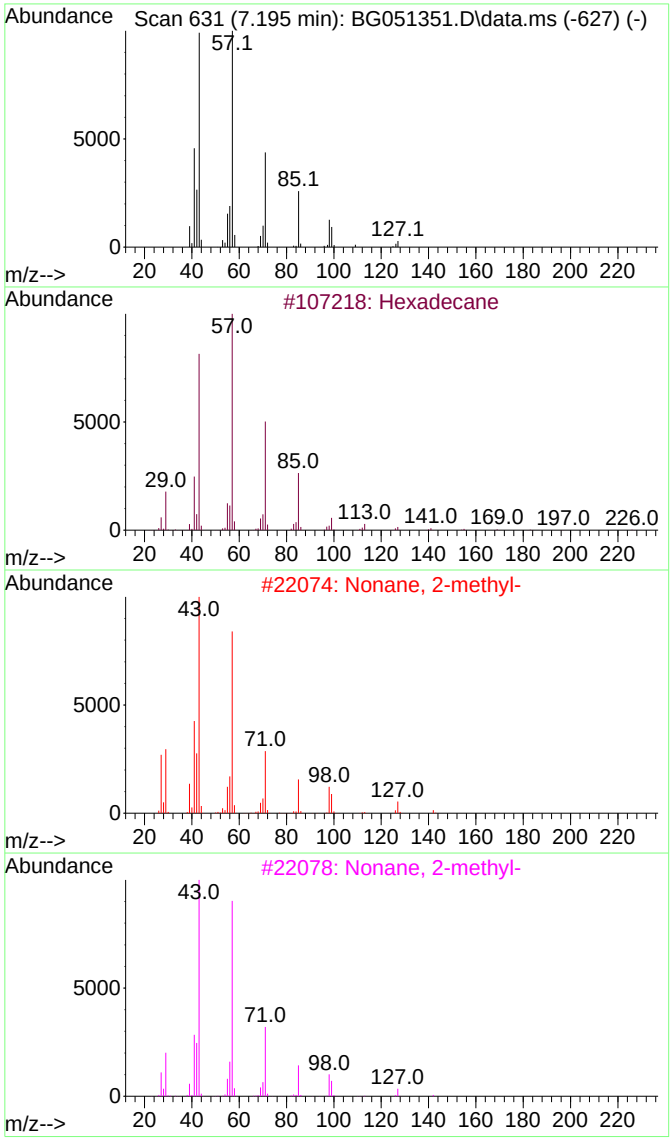
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

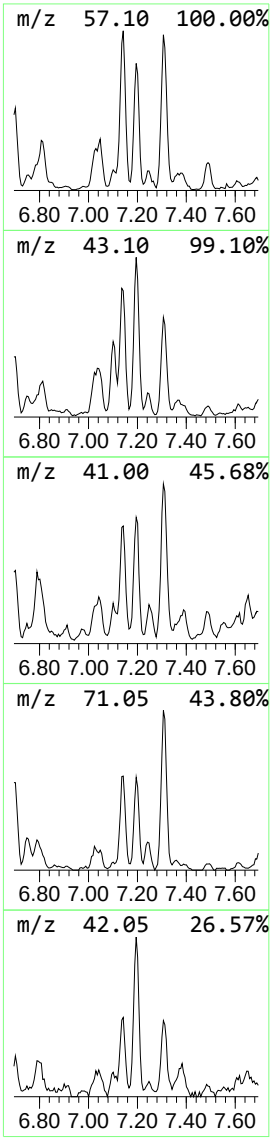
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.195	18.30 ng/ul	177414	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	72
2	Nonane, 2-methyl-	142	C10H22	000871-83-0	64
3	Nonane, 2-methyl-	142	C10H22	000871-83-0	64
4	Decane	142	C10H22	000124-18-5	64
5	Sulfurous acid, hexyl octyl ester	278	C14H30O3S	1000309-13-0	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

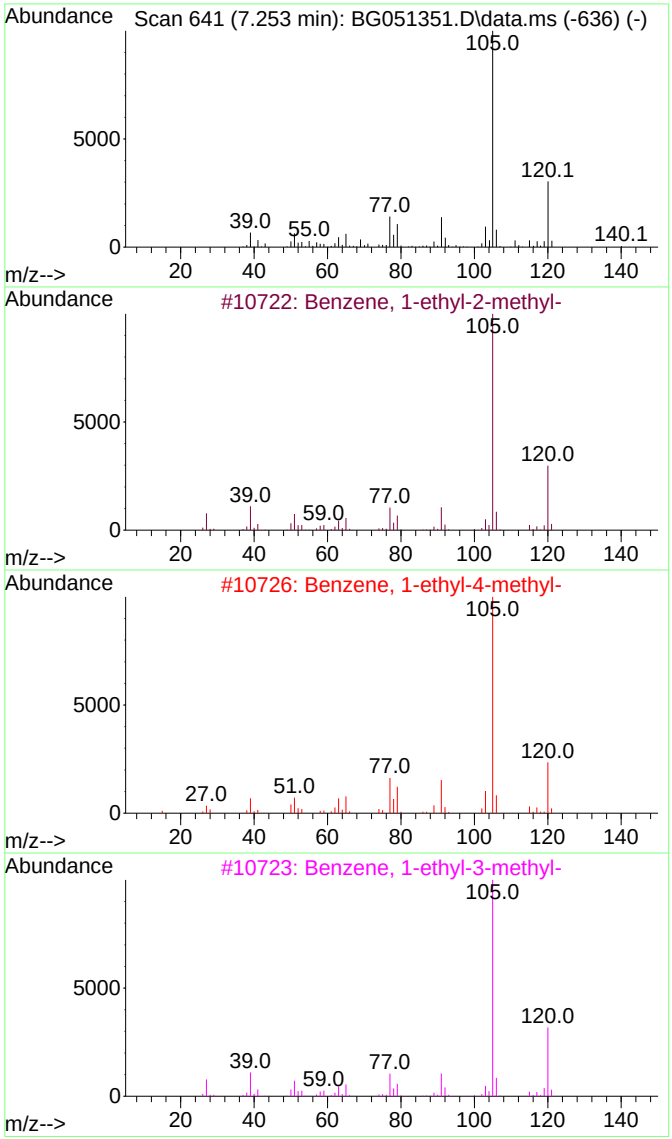
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

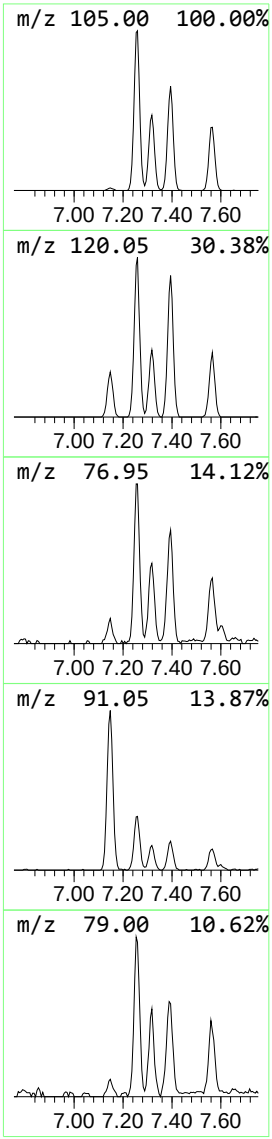
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.253	40.20 ng/ul	389817	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

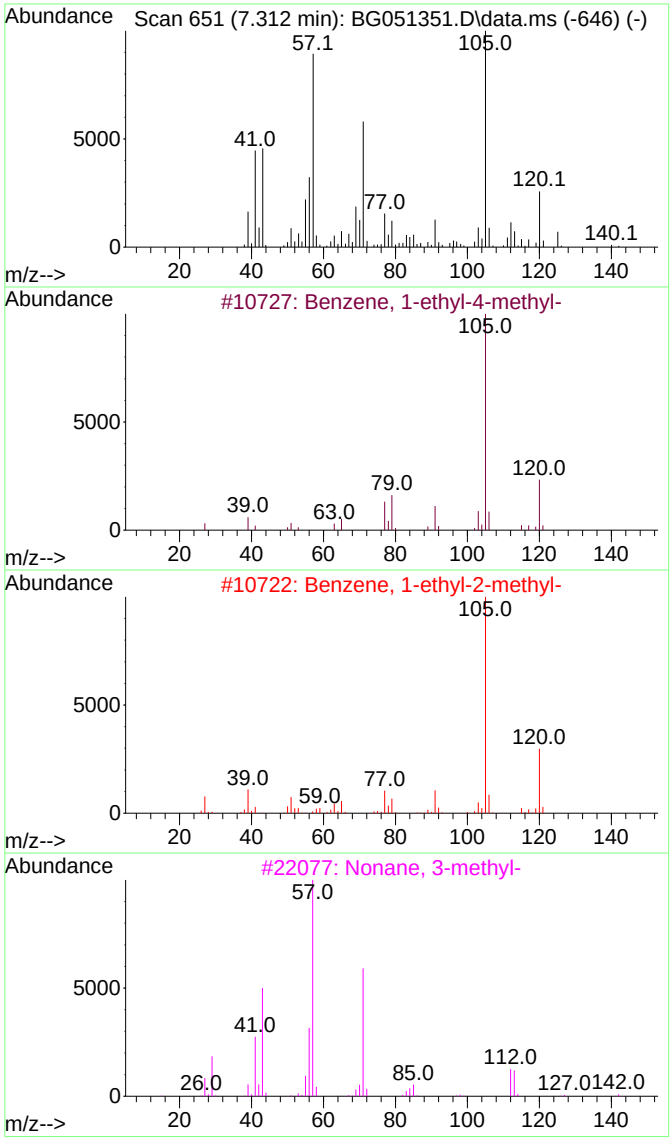
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

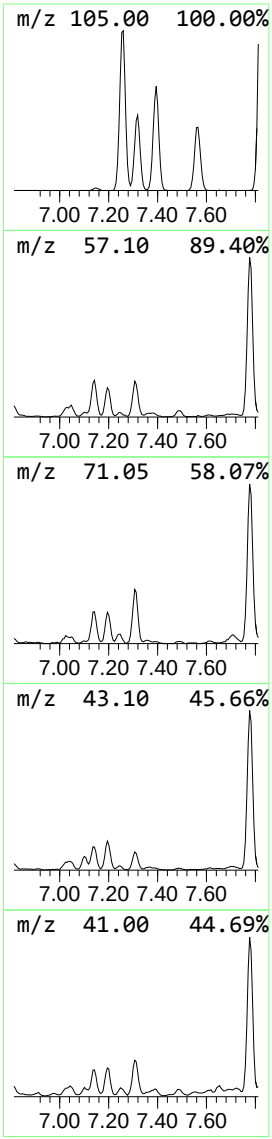
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-4-methyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.312	42.27 ng/ul	409855	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	70
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	60
3	Nonane, 3-methyl-	142	C10H22	005911-04-6	60
4	Nonane, 3-methyl-	142	C10H22	005911-04-6	53
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	52





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

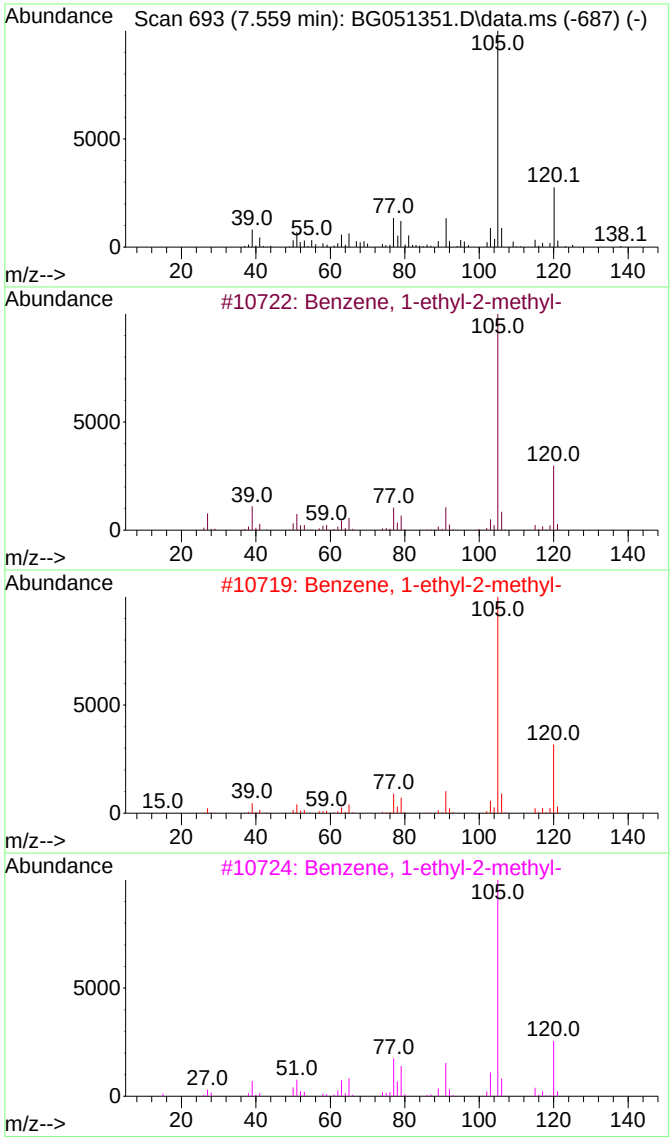
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

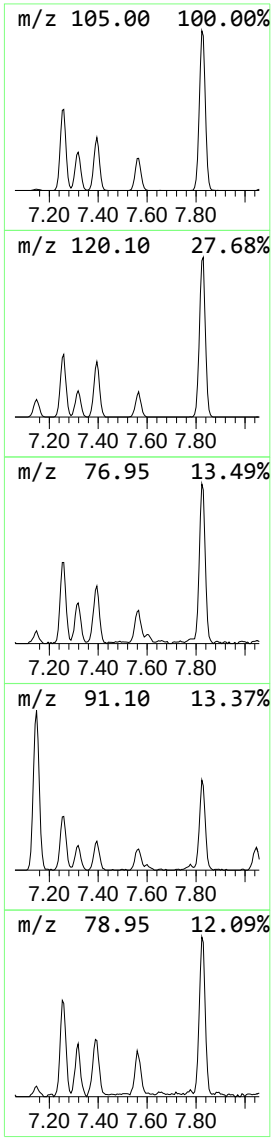
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.559	18.86 ng/ul	182869	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

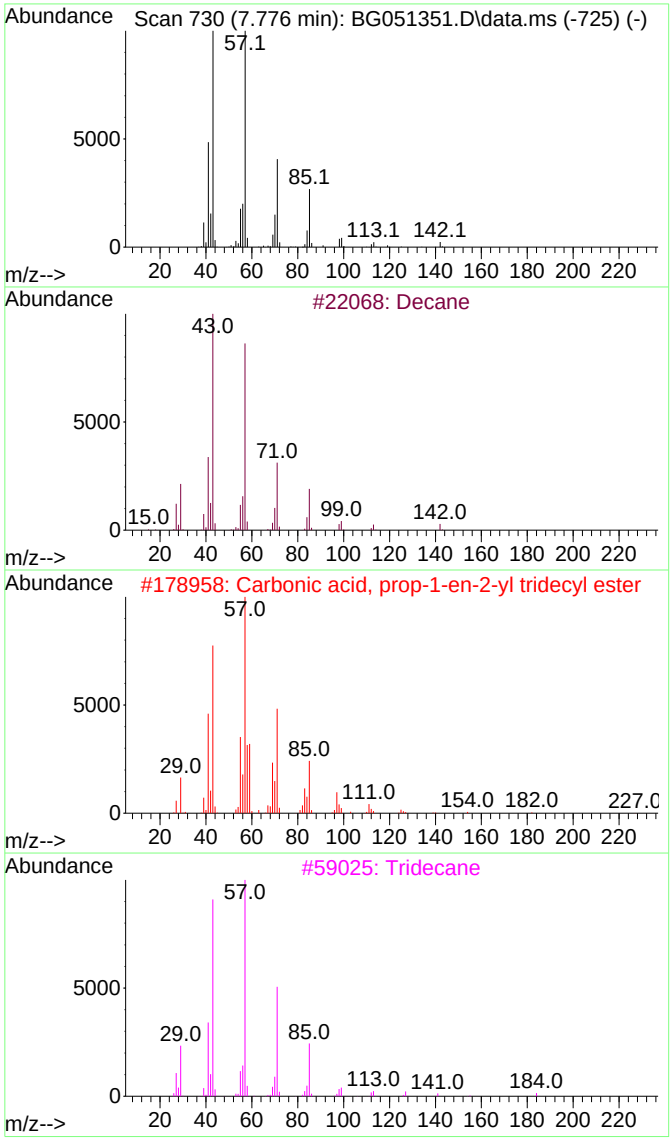
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

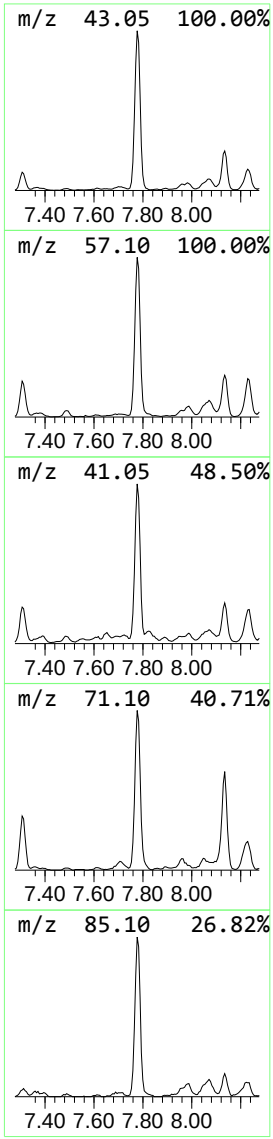
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.776	97.87 ng/ul	949007	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	91
2	Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	83
3	Tridecane	184	C13H28	000629-50-5	72
4	Decane	142	C10H22	000124-18-5	72
5	Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

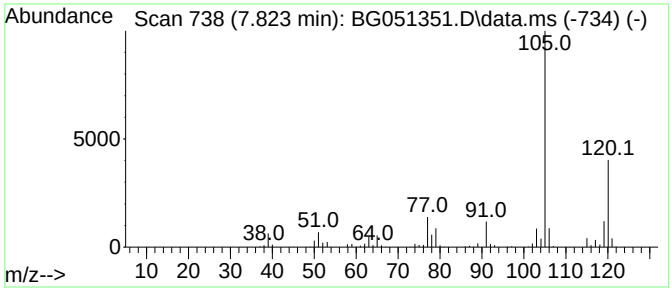
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

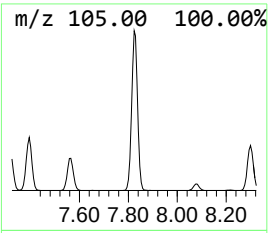
****

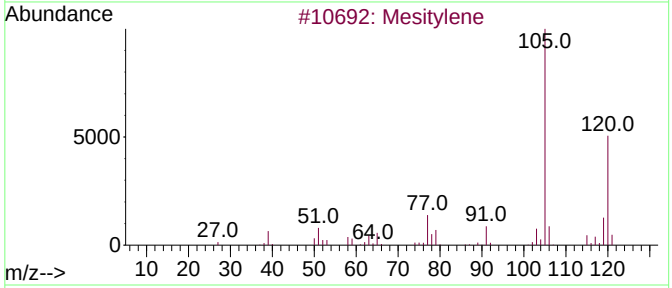
Peak Number 11 Mesitylene Concentration Rank 5

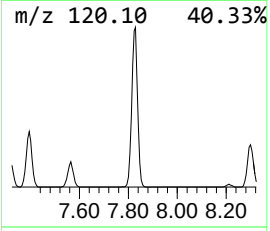
R.T.	EstConc	Area	Relative to ISTD	R.T.
7.823	77.83 ng/ul	754702	1,4-Dichlorobenzene-d4	8.194

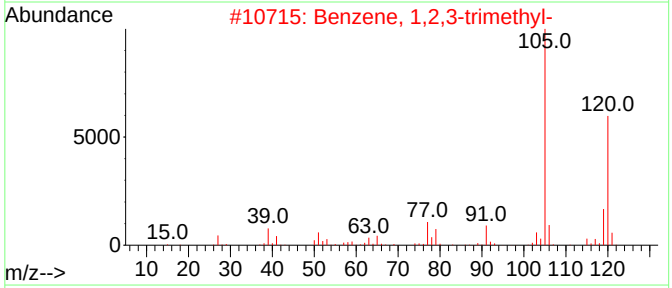
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5	Mesitylene	120	C9H12	000108-67-8	91

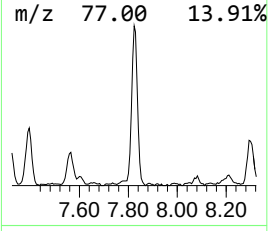


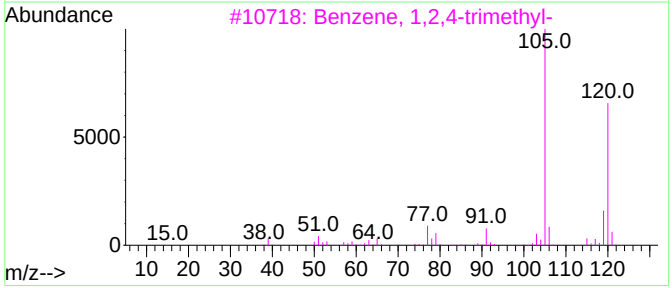


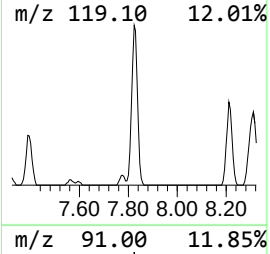


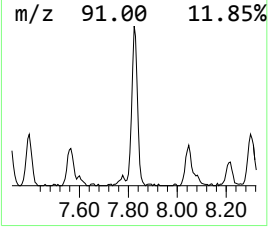












Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

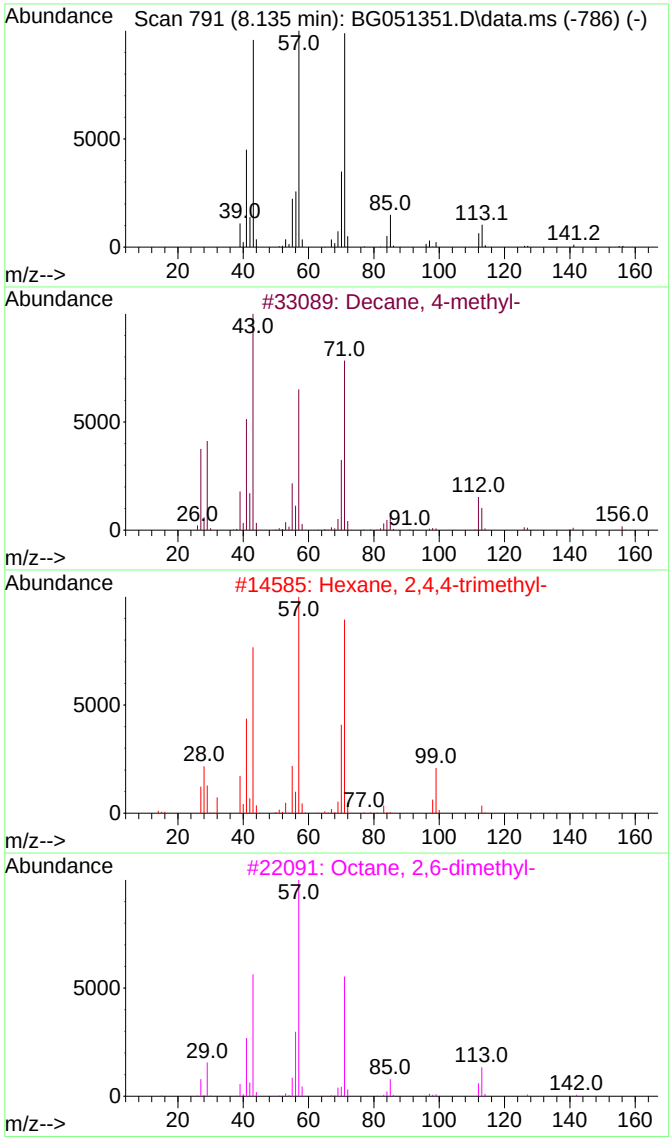
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

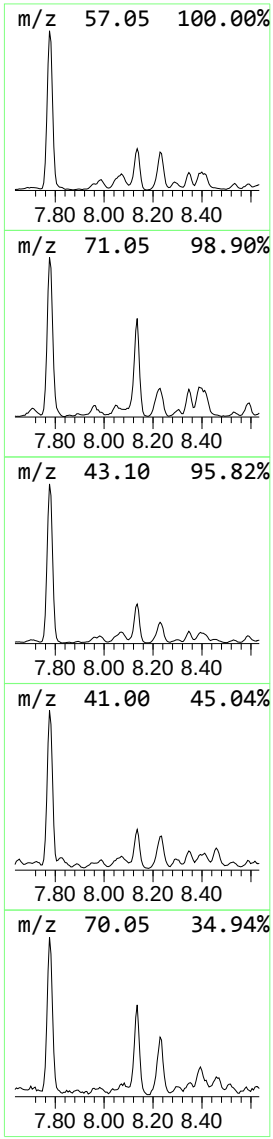
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.135	28.56 ng/ul	276956	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-	156	C11H24	002847-72-5	72
2	Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	64
3	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	64
5	Sulfurous acid, octyl 2-propyl e...	236	C11H24O3S	1000309-11-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

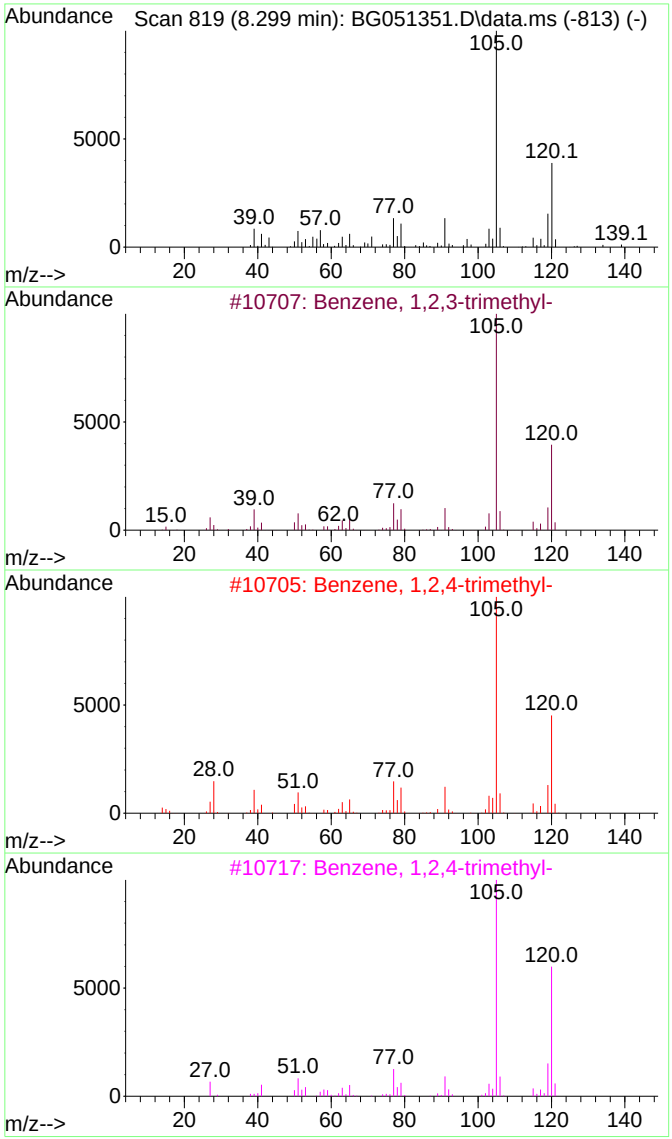
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

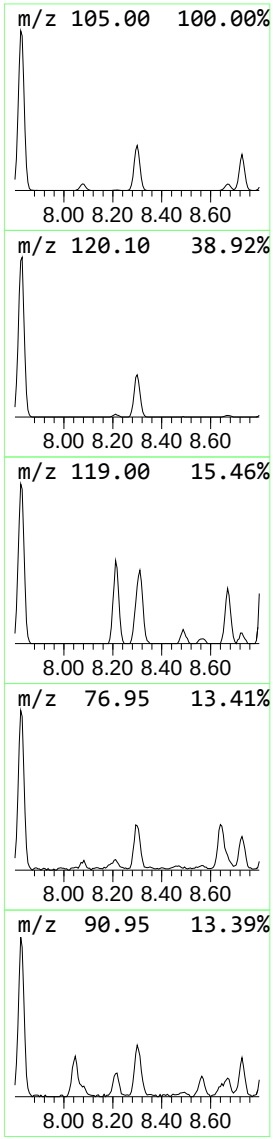
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1,2,3-trimethyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.299	28.95 ng/ul	280704	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	93
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

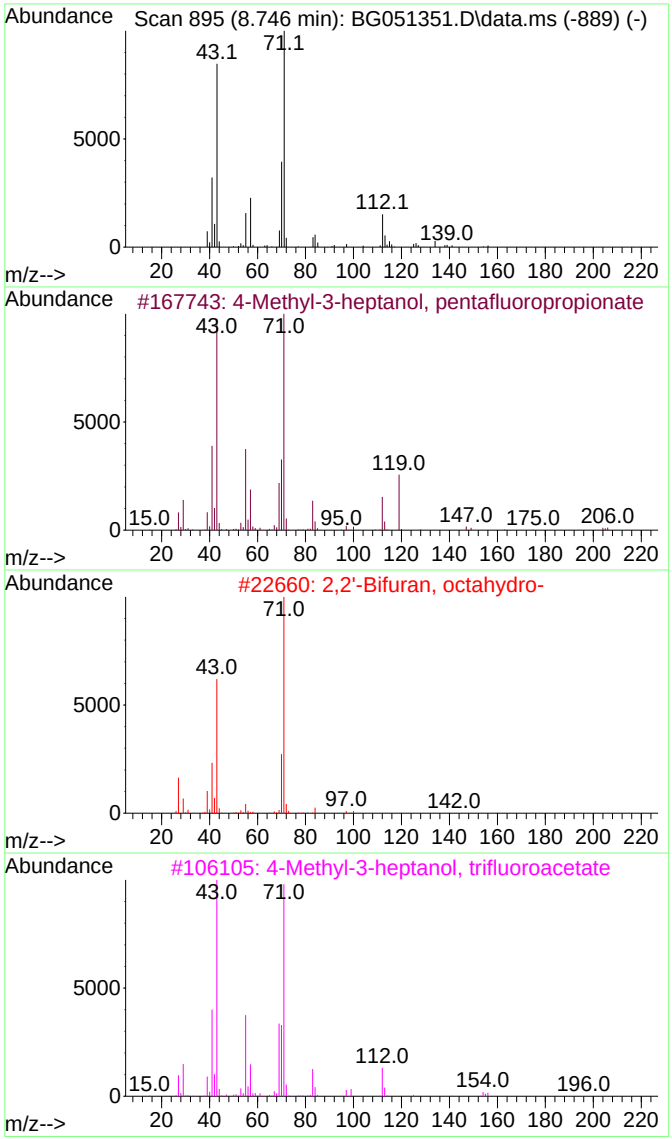
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

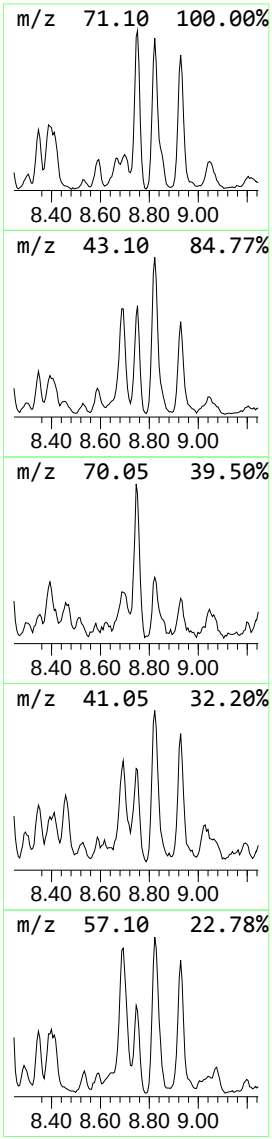
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 4-Methyl-3-heptanol, pentaf... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.746	41.34 ng/ul	400854	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	78
2	2,2'-Bifuran, octahydro-	142	C8H14O2	001592-33-2	72
3	4-Methyl-3-heptanol, trifluoroac...	226	C10H17F3O2	1000365-51-8	56
4	Tridecane, 4-methyl-	198	C14H30	026730-12-1	56
5	Pentane, 3-bromo-	150	C5H11Br	001809-10-5	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

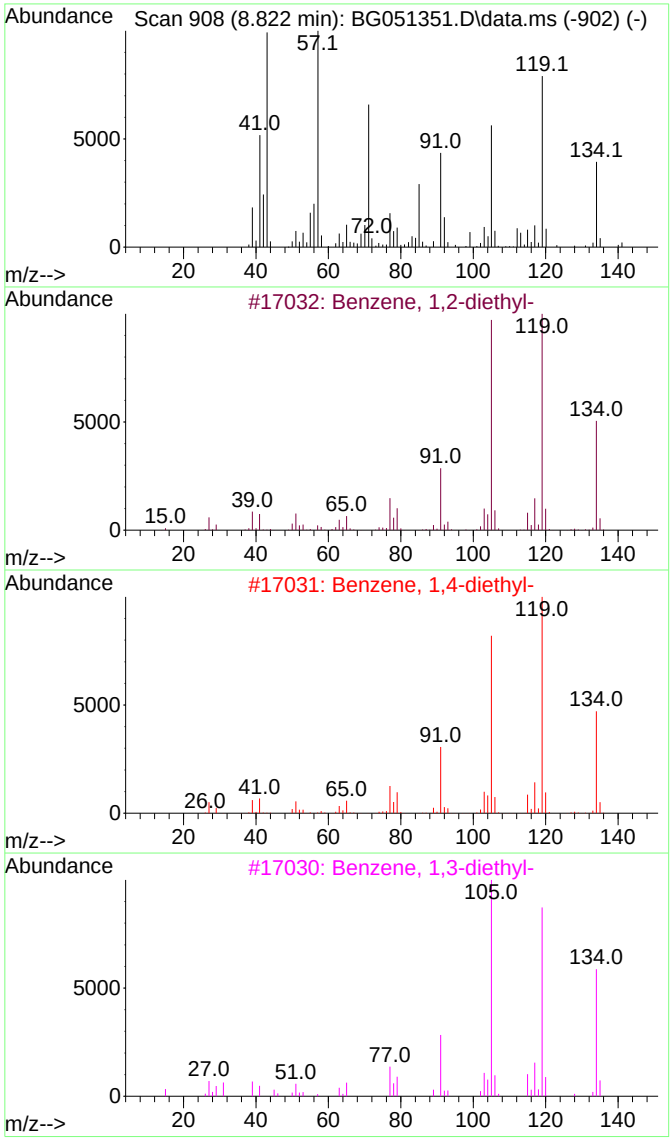
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

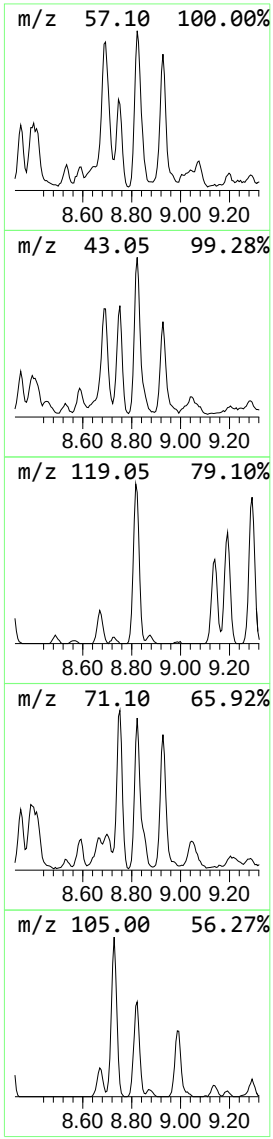
****

Peak Number 15 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.822	63.95 ng/ul	620115	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
2	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	70
4	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	70
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

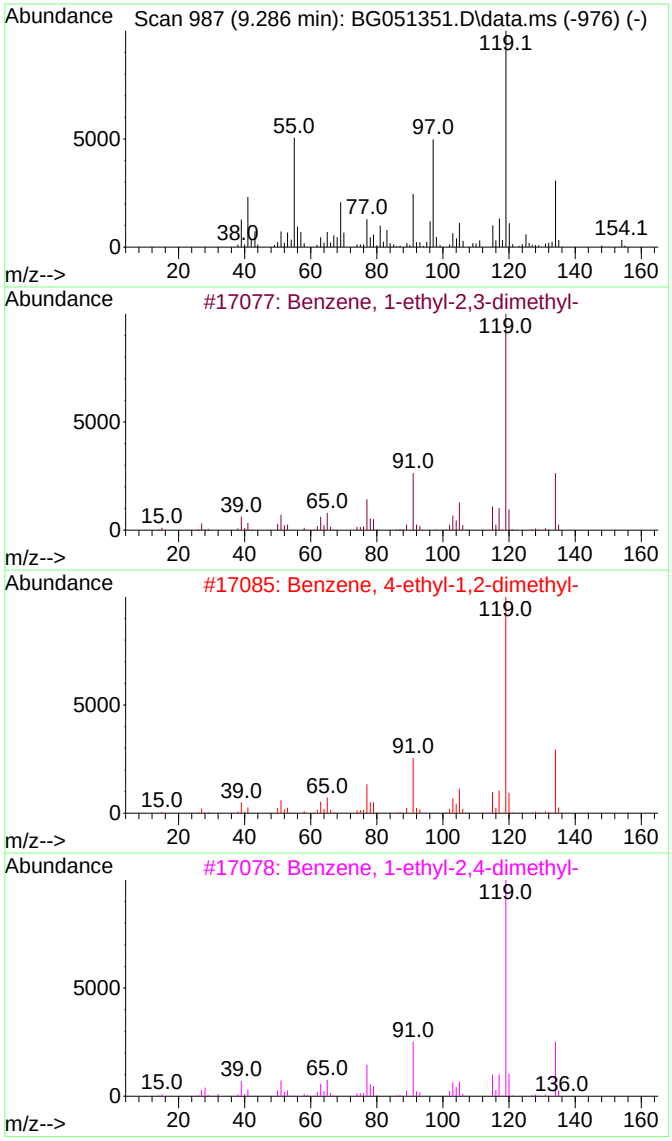
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

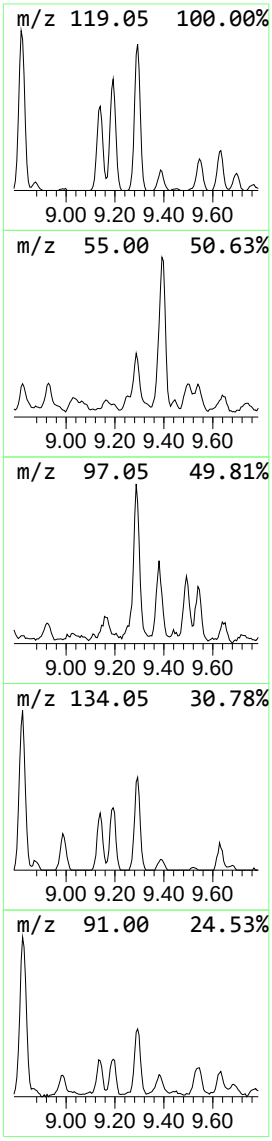
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.286	36.51 ng/ul	354034	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93
5	o-Cymene	134	C10H14	000527-84-4	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

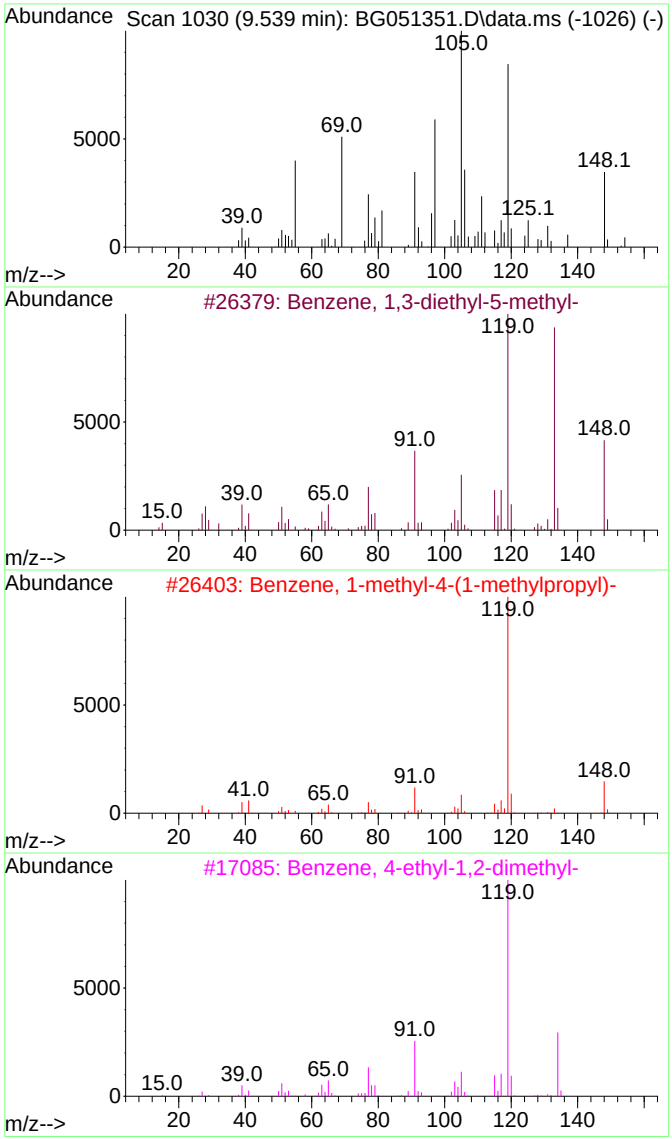
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

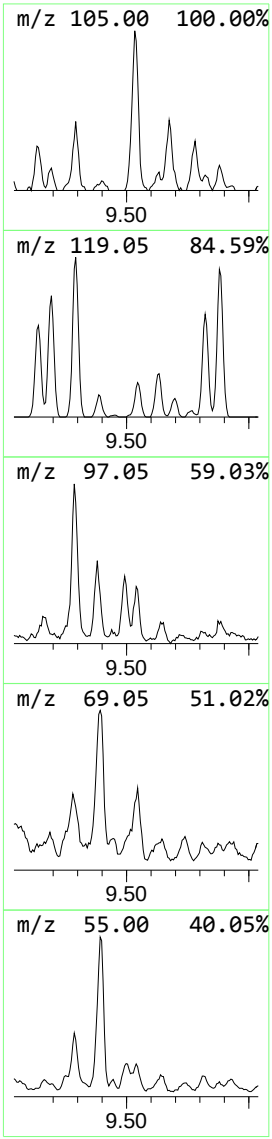
****

Peak Number 17 unknown-02 Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.539	19.92 ng/ul	193133	1,4-Dichlorobenzene-d4	8.194

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	46
2	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	38
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	35
4	3,4-Dihydro-2-[1H]quinoxalinone	148	C8H8N2O	059564-59-9	30
5	4-Methylphthalaldehyde	148	C9H8O2	015158-36-8	27





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

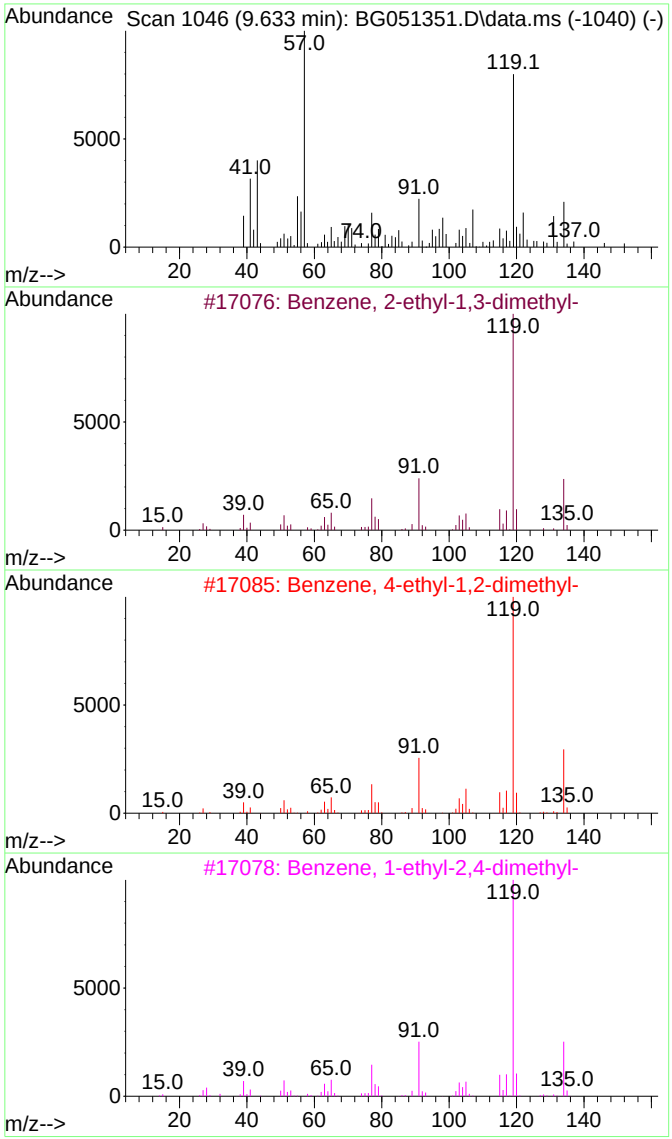
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

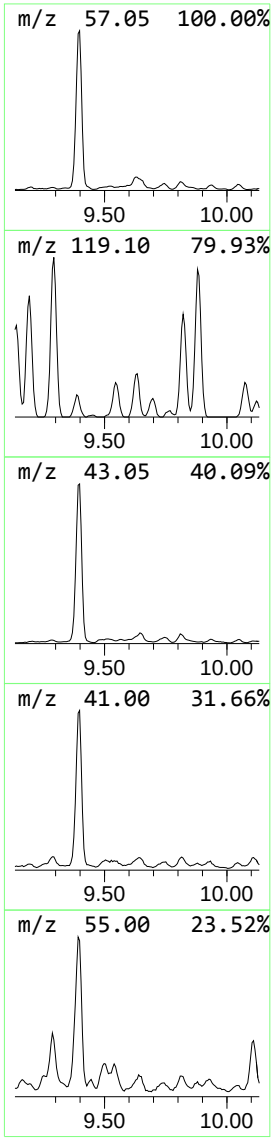
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.633	18.27 ng/ul	267399	Naphthalene-d8	11.025

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91
2	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	90
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

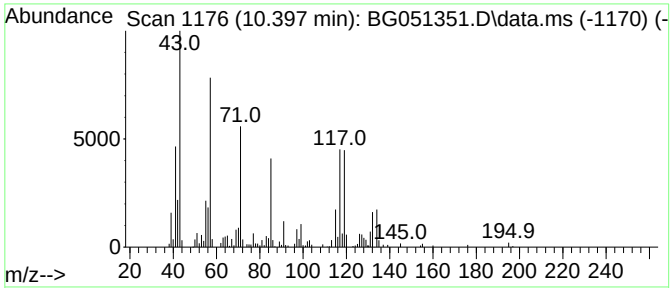
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

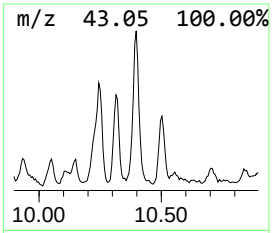
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

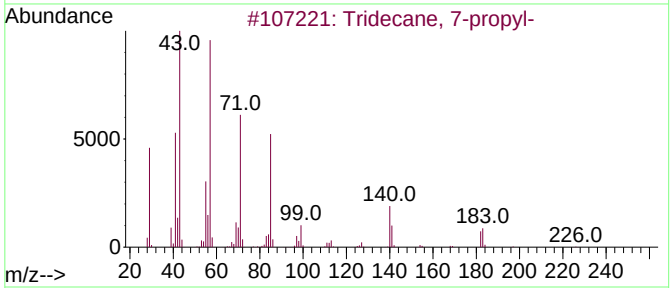
****
Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 45

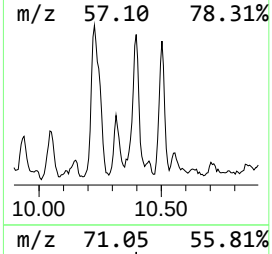
R.T.	EstConc	Area	Relative to ISTD	R.T.
10.397	19.95 ng/ul	291927	Naphthalene-d8	11.025

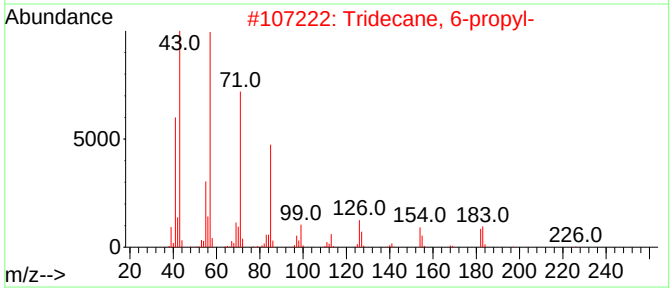
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 7-propyl-	226	C16H34	055045-09-5	38
2	Tridecane, 6-propyl-	226	C16H34	055045-10-8	38
3	Heptadecane, 8-methyl-	254	C18H38	013287-23-5	38
4	Tetracosane	338	C24H50	000646-31-1	38
5	Dodecane, 3-methyl-	184	C13H28	017312-57-1	38

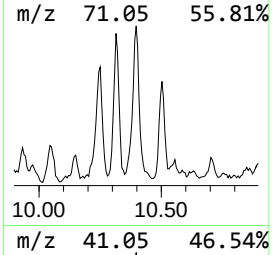


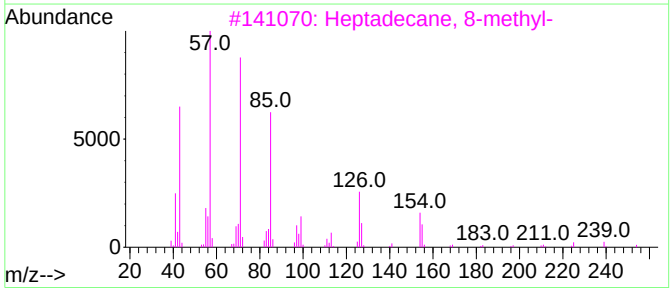


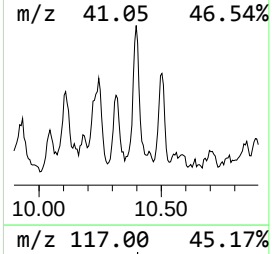


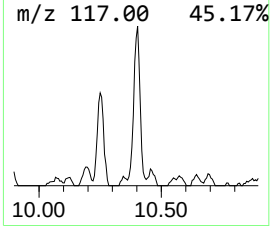












Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

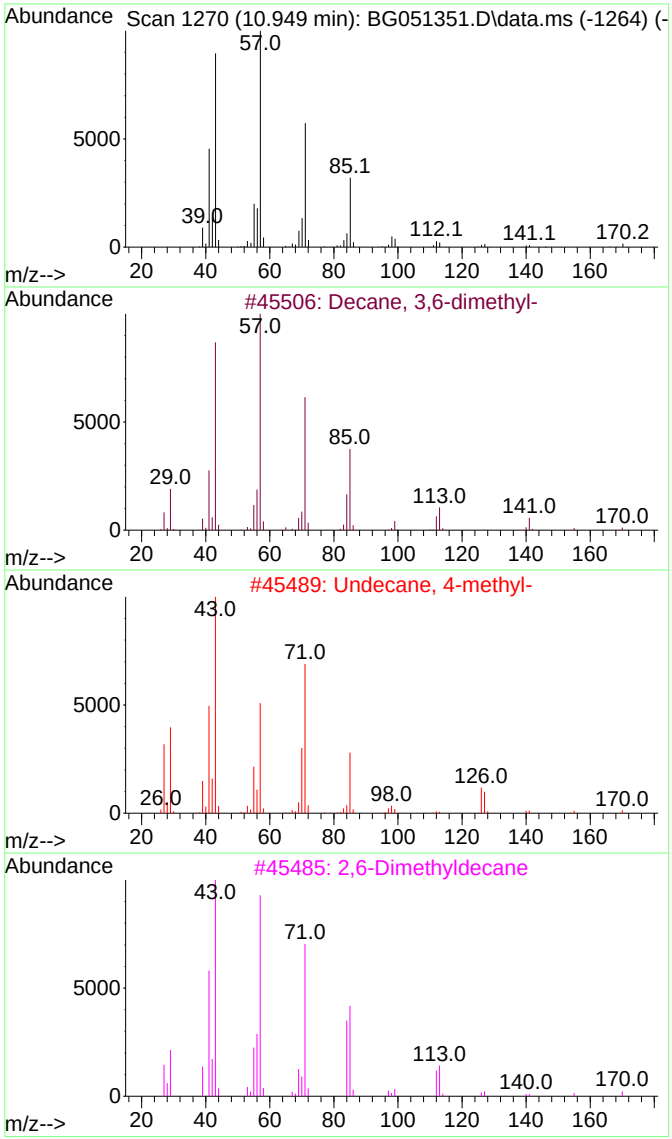
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

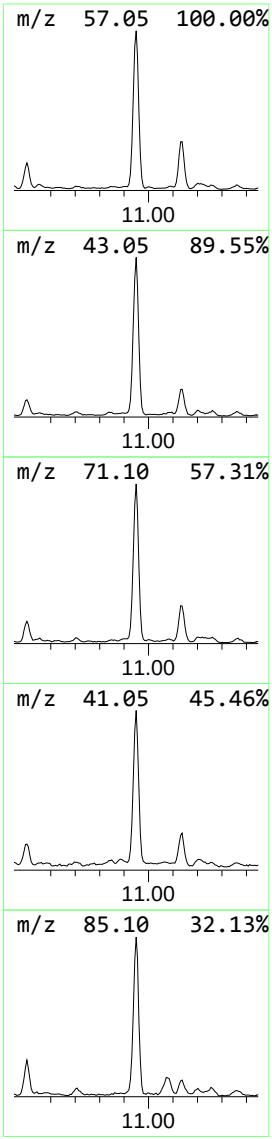
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.949	60.44 ng/ul	884348	Naphthalene-d8	11.025

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	86
2	Undecane, 4-methyl-	170	C12H26	002980-69-0	72
3	2,6-Dimethyldecane	170	C12H26	013150-81-7	64
4	Nonane	128	C9H20	000111-84-2	59
5	Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

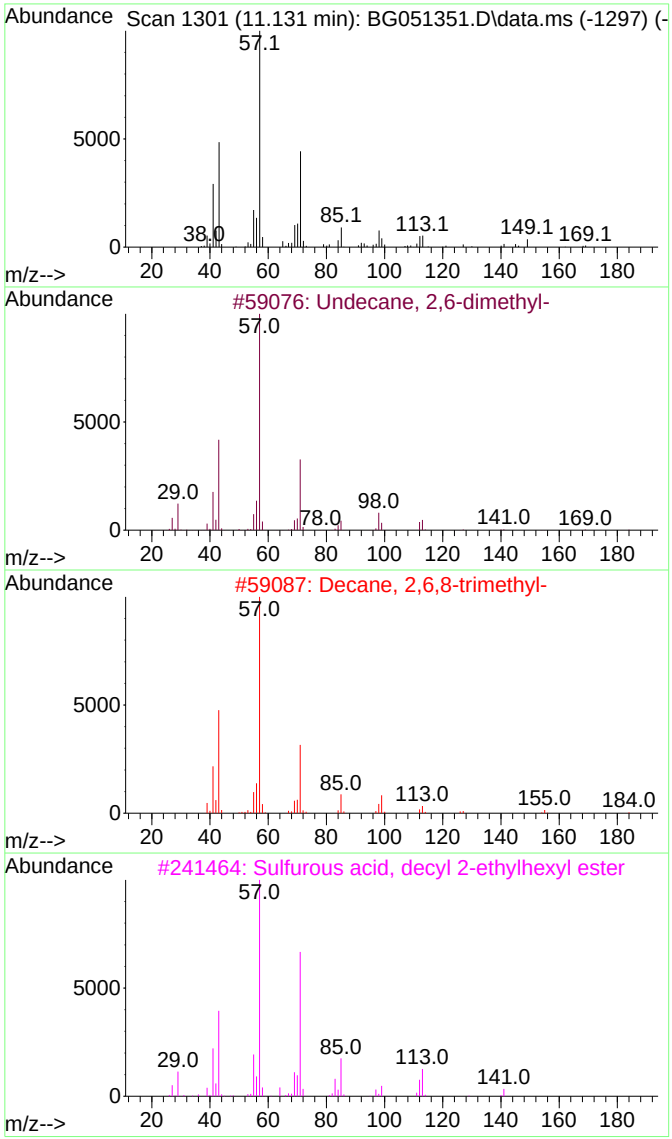
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

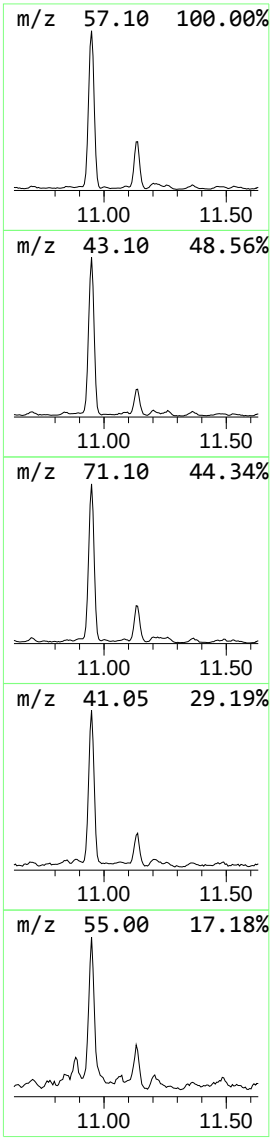
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.131	15.73 ng/ul	230143	Naphthalene-d8	11.025

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	86
2	Decane, 2,6,8-trimethyl-	184	C13H28	062108-26-3	78
3	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	78
4	Undecane, 3,6-dimethyl-	184	C13H28	017301-28-9	78
5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

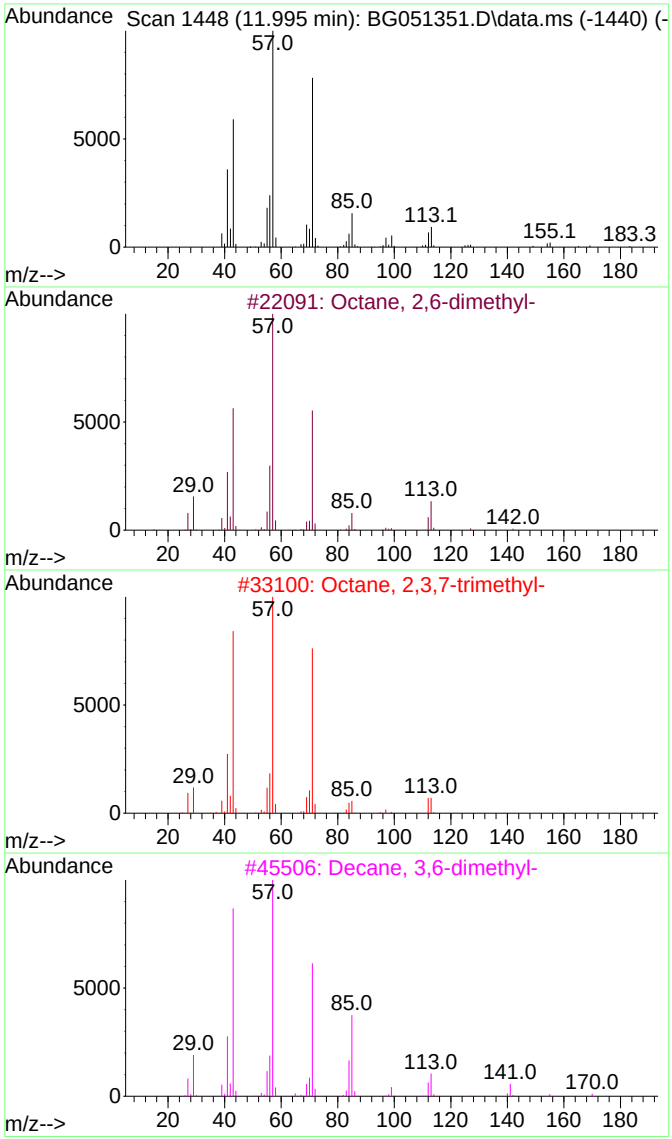
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

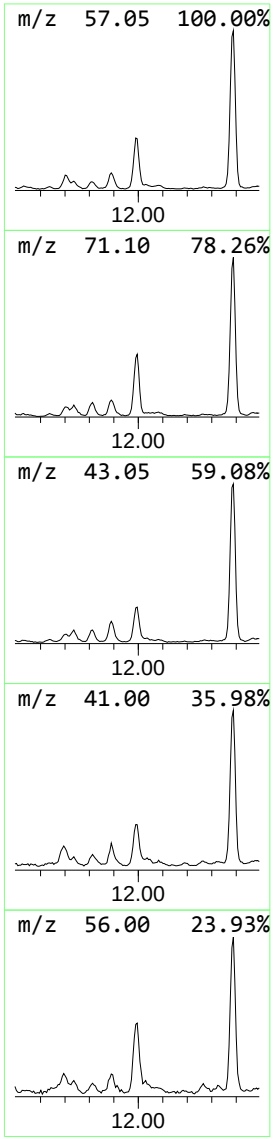
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.995	17.56 ng/ul	256978	Naphthalene-d8	11.025

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
2	Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	72
3	Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	64
4	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	64
5	Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

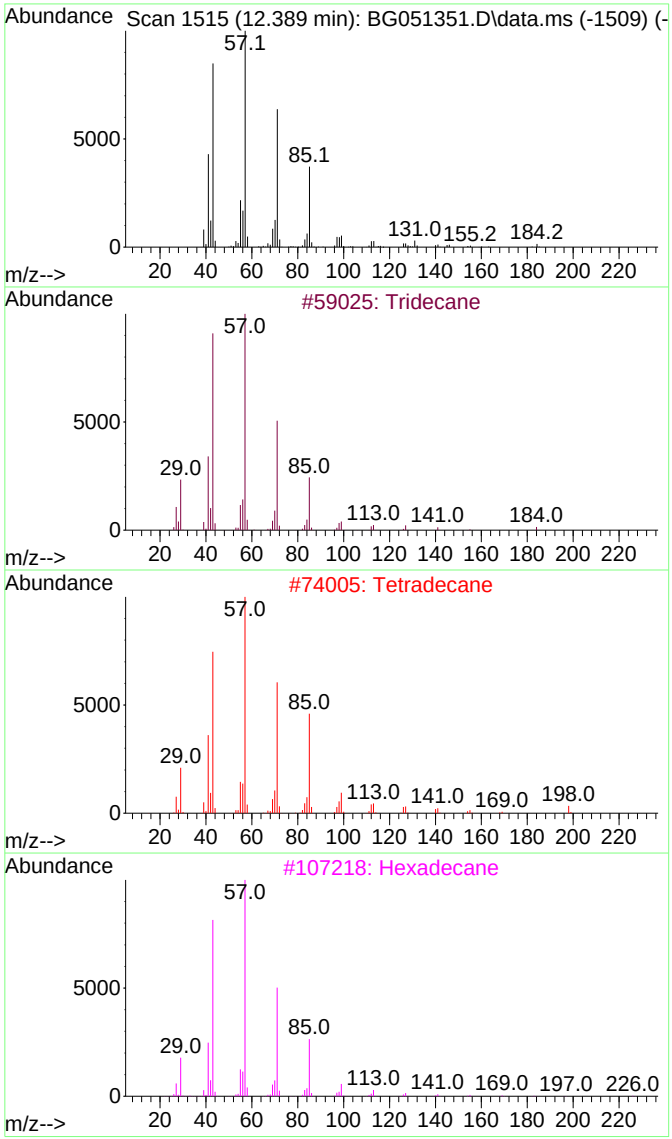
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

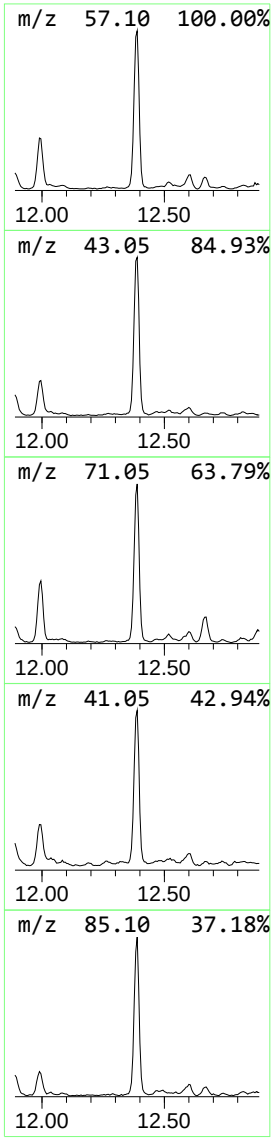
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.389	53.15 ng/ul	777637	Naphthalene-d8	11.025

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane	184	C13H28	000629-50-5	87
2	Tetradecane	198	C14H30	000629-59-4	86
3	Hexadecane	226	C16H34	000544-76-3	86
4	Undecane, 6-ethyl-	184	C13H28	017312-60-6	86
5	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

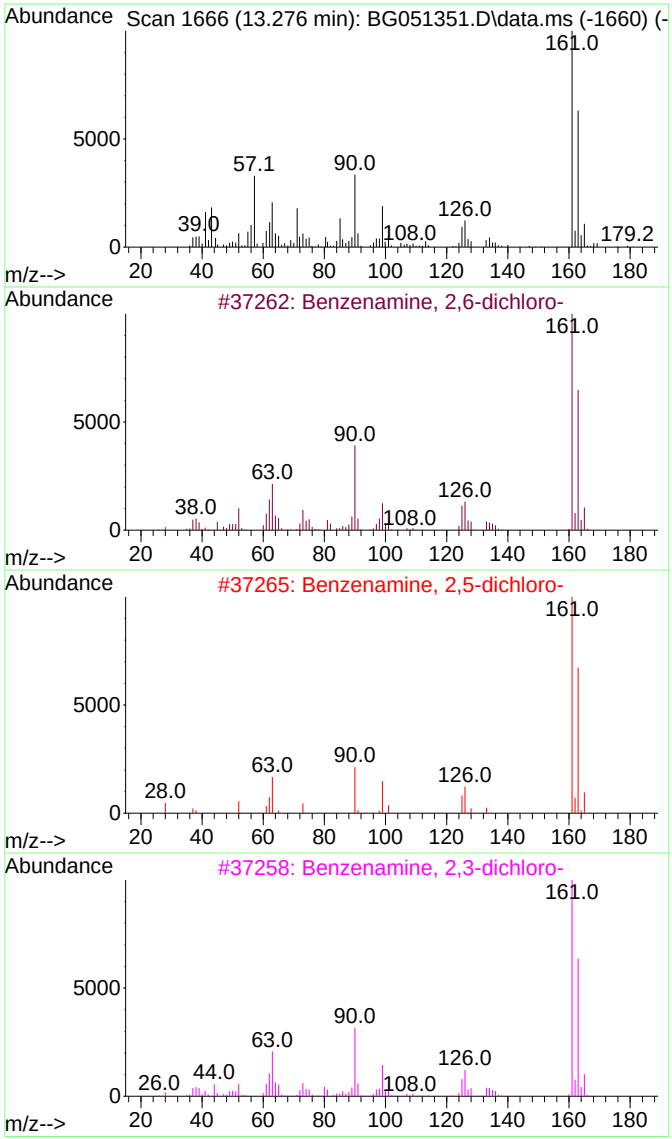
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

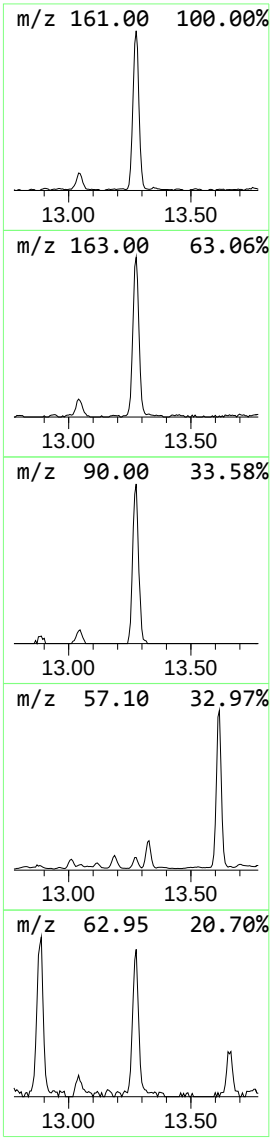
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Benzenamine, 2,6-dichloro- Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.276	18.61 ng/ul	316992	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	97
2	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	97
3	Benzenamine, 2,3-dichloro-	161	C6H5Cl2N	000608-27-5	96
4	Benzenamine, 2,6-dichloro-	161	C6H5Cl2N	000608-31-1	95
5	Benzenamine, 2,5-dichloro-	161	C6H5Cl2N	000095-82-9	95





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

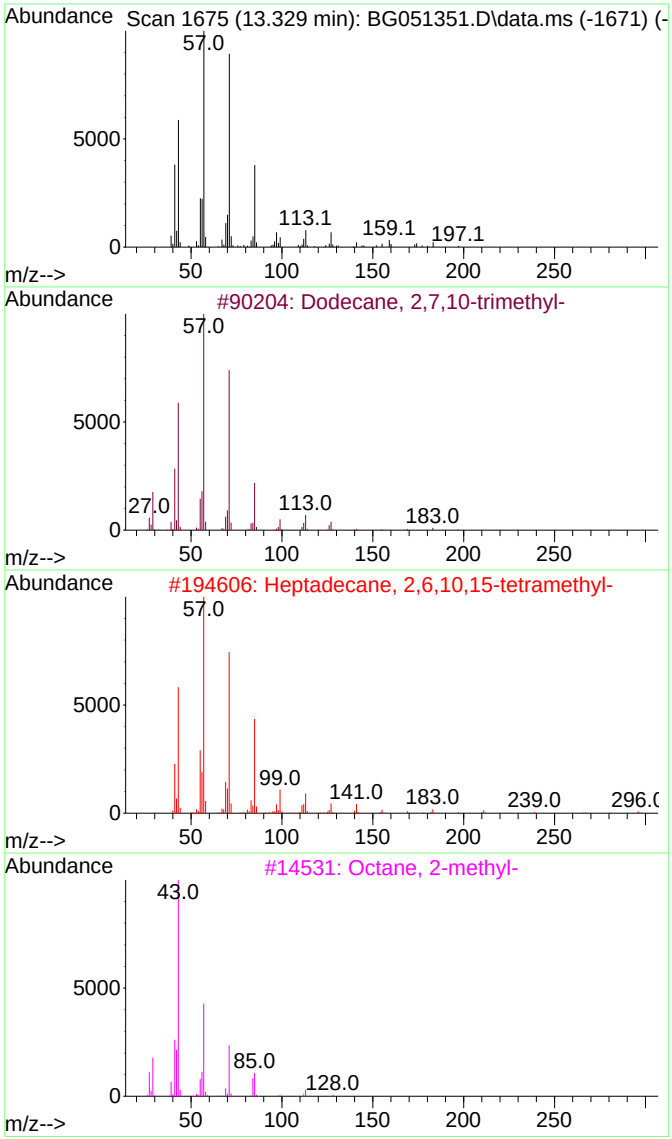
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

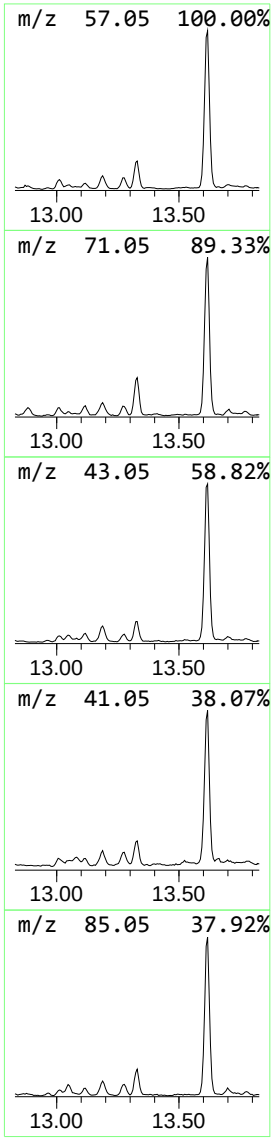
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 72

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.329	11.01 ng/ul	187573	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	90
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	83
3	Octane, 2-methyl-	128	C9H20	003221-61-2	80
4	Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	72
5	Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	59





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

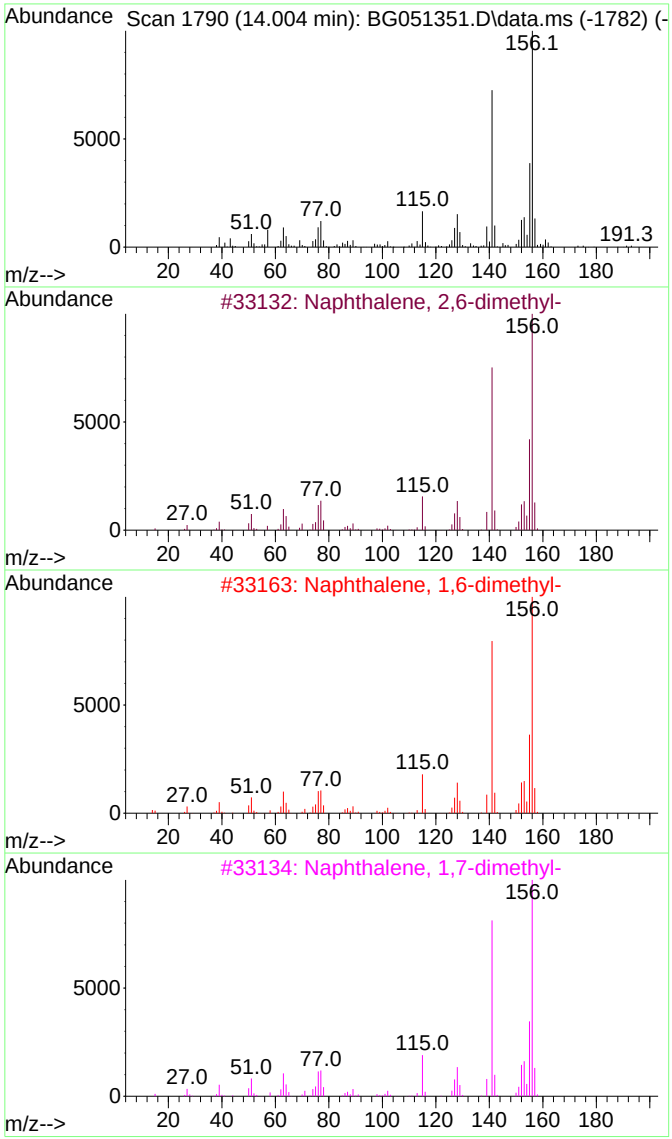
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

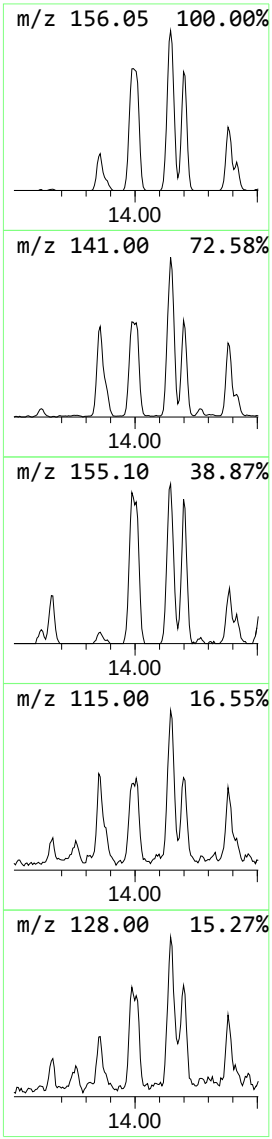
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 Naphthalene, 2,6-dimethyl- Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.004	23.91 ng/ul	407212	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

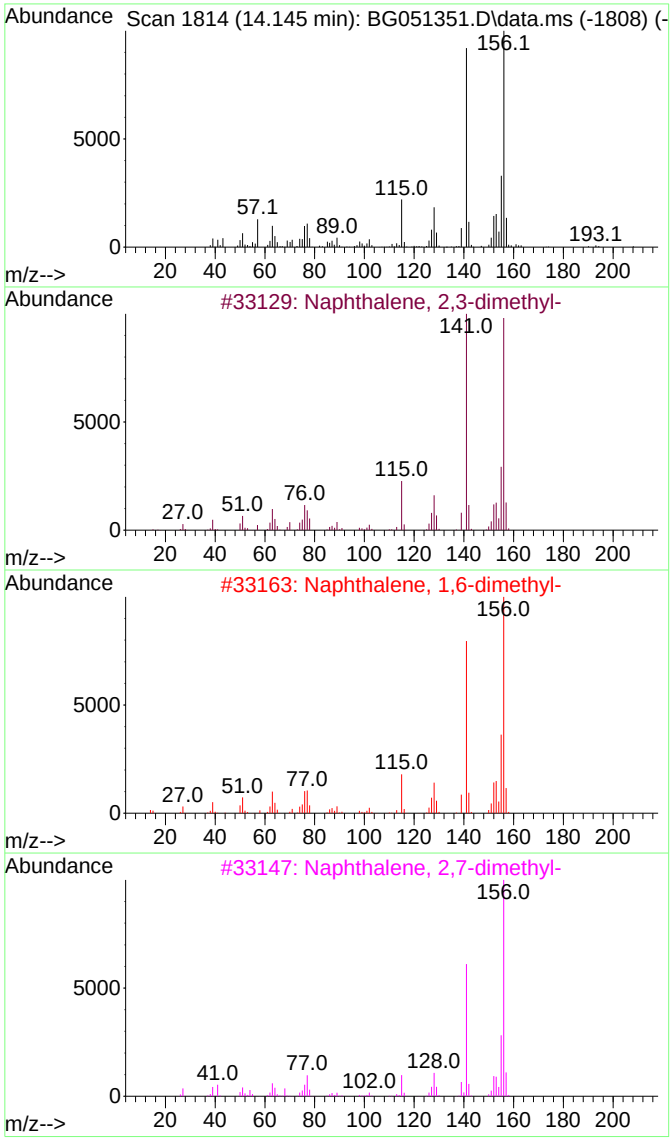
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

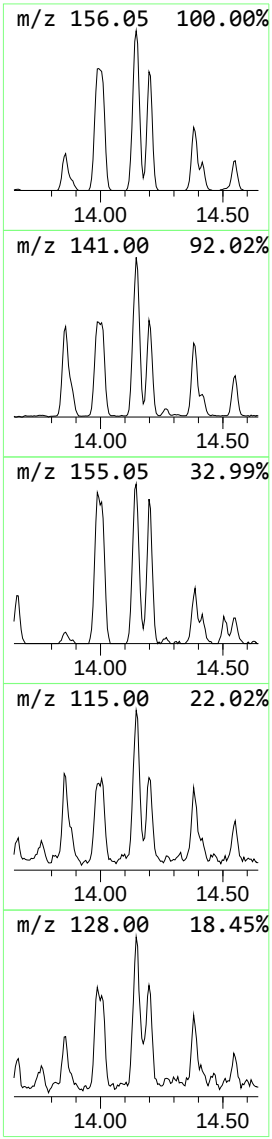
****

Peak Number 30 Naphthalene, 2,3-dimethyl- Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.145	28.45 ng/ul	484553	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
4	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

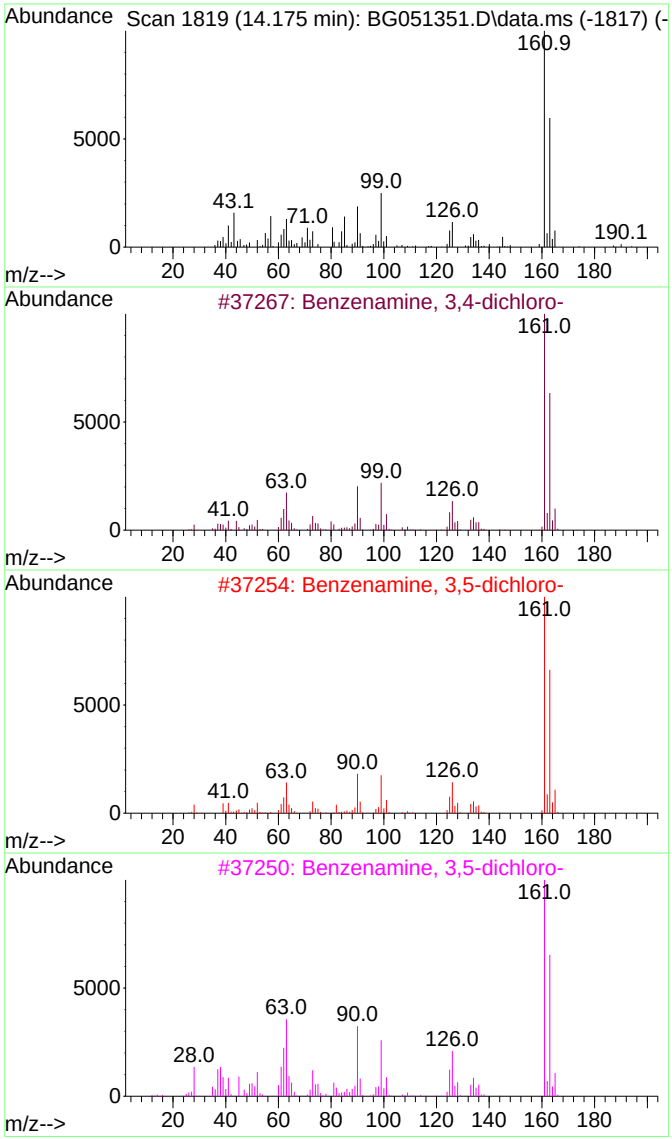
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

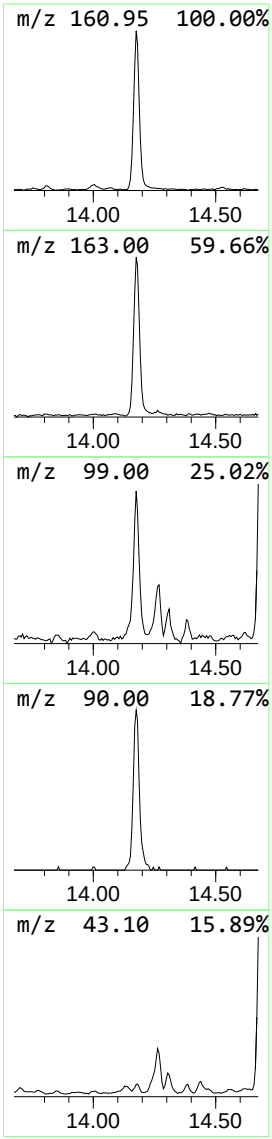
****

Peak Number 31 Benzenamine, 3,4-dichloro- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.175	31.10 ng/ul	529639	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 3,4-dichloro-	161	C6H5Cl2N	000095-76-1	96
2	Benzenamine, 3,5-dichloro-	161	C6H5Cl2N	000626-43-7	95
3	Benzenamine, 3,5-dichloro-	161	C6H5Cl2N	000626-43-7	95
4	Benzenamine, 3,5-dichloro-	161	C6H5Cl2N	000626-43-7	94
5	Benzenamine, 2,4-dichloro-	161	C6H5Cl2N	000554-00-7	94





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

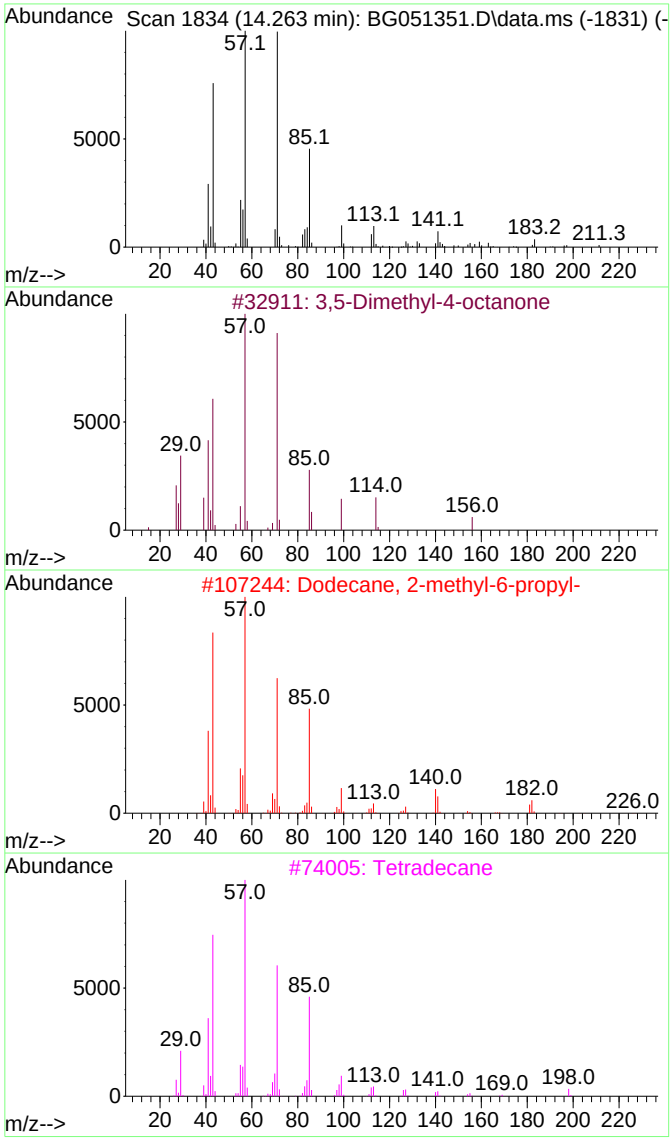
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

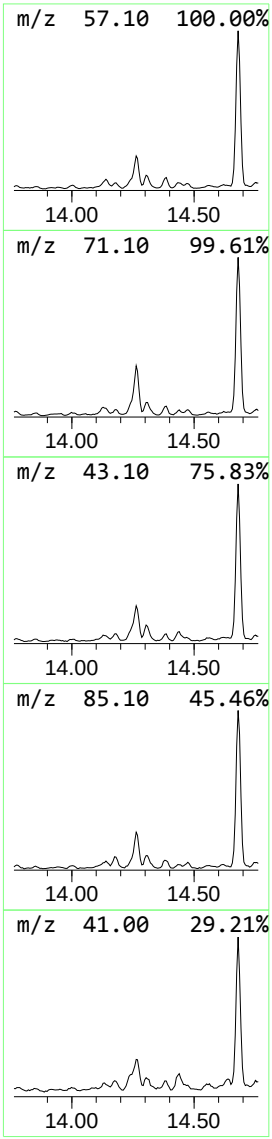
****

Peak Number 32 3,5-Dimethyl-4-octanone Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.263	18.18 ng/ul	309665	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethyl-4-octanone	156	C10H20O	007335-17-3	53
2	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	53
3	Tetradecane	198	C14H30	000629-59-4	53
4	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	53
5	Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

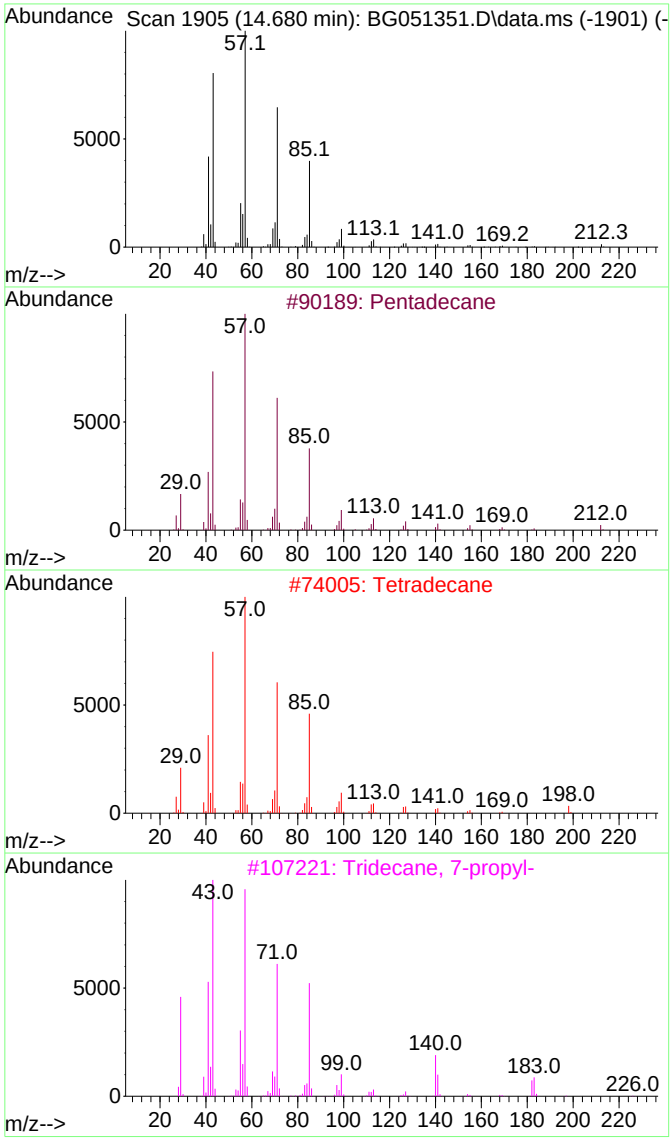
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

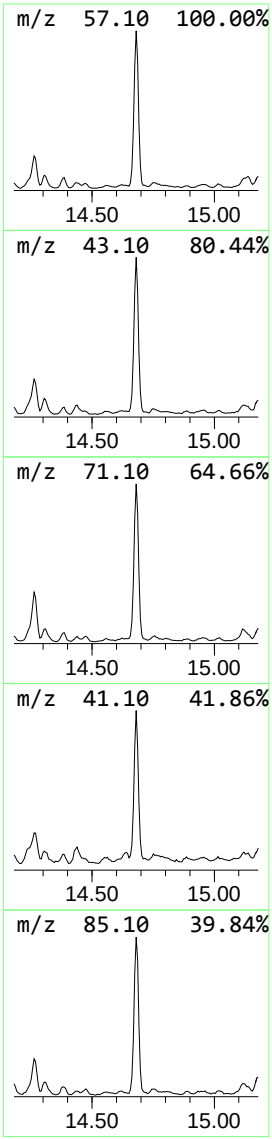
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 35 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.680	53.11 ng/ul	904495	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecane	212	C15H32	000629-62-9	91
2	Tetradecane	198	C14H30	000629-59-4	90
3	Tridecane, 7-propyl-	226	C16H34	055045-09-5	86
4	Octacosane	394	C28H58	000630-02-4	78
5	Tetradecane	198	C14H30	000629-59-4	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

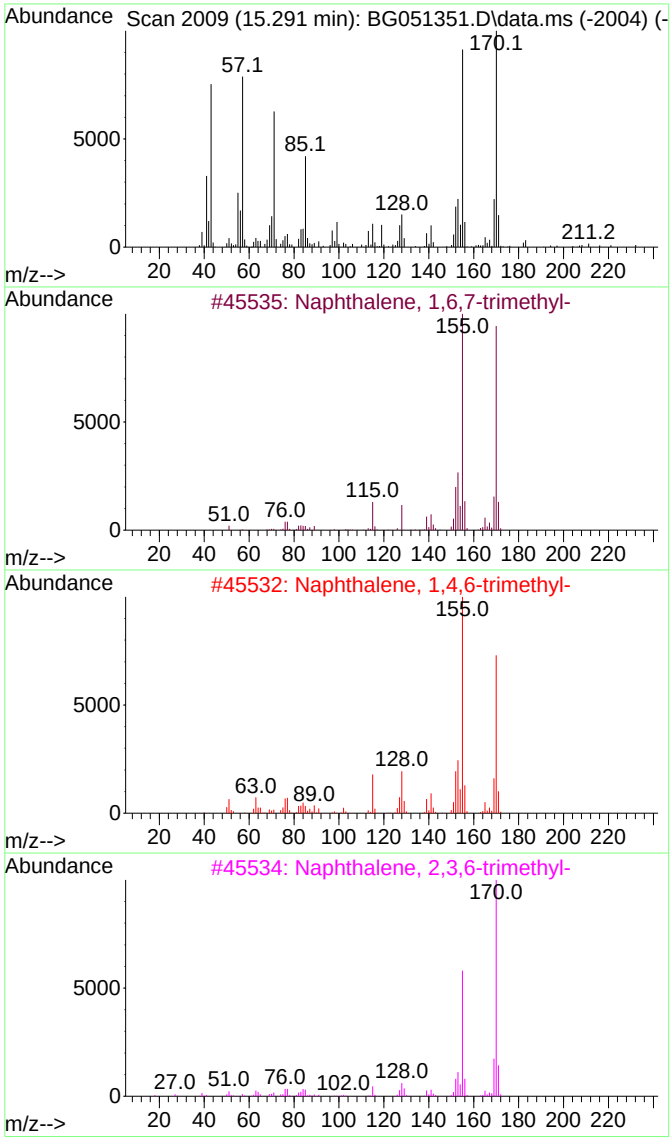
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

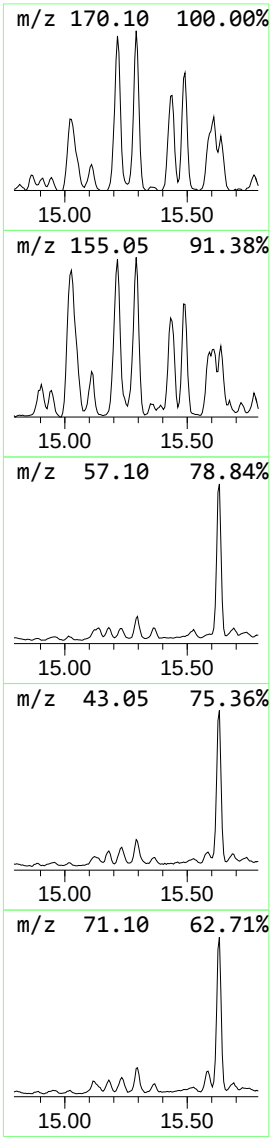
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 Naphthalene, 1,6,7-trimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.291	20.19 ng/ul	343844	Acenaphthene-d10	14.827

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

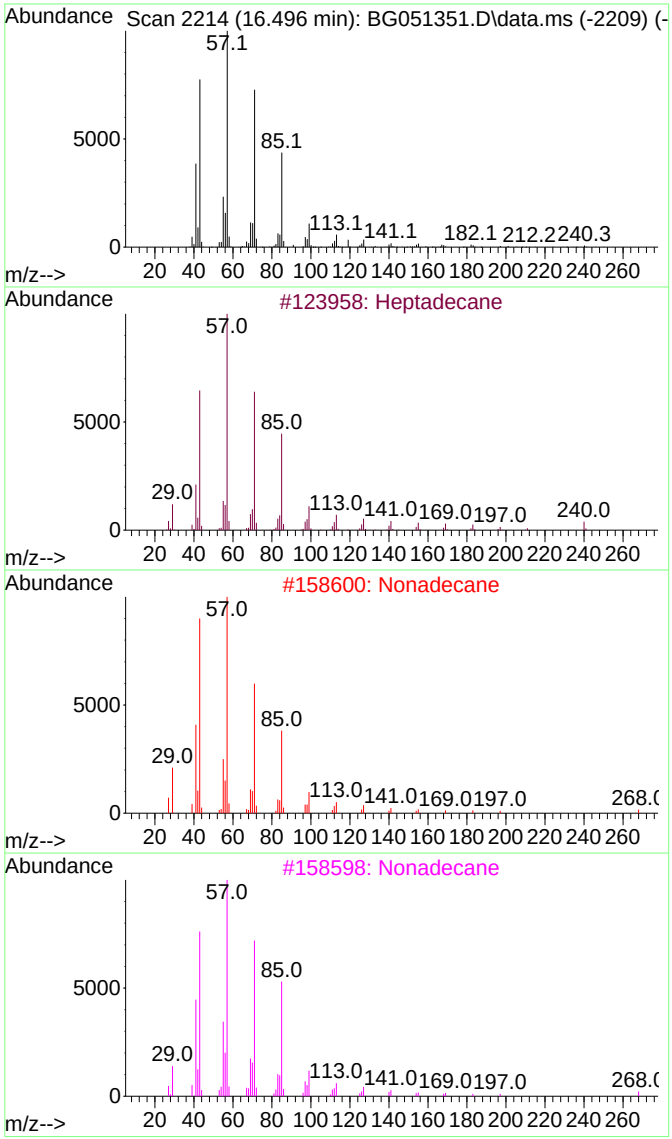
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

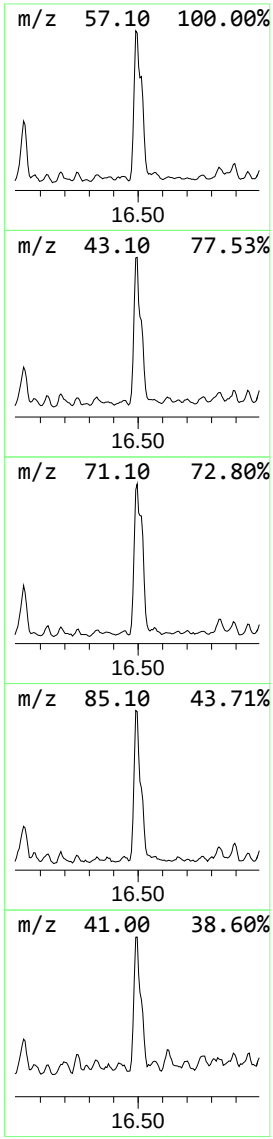
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.496	78.26 ng/ul	1550930	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Nonadecane	268	C19H40	000629-92-5	91
3	Nonadecane	268	C19H40	000629-92-5	91
4	Heneicosane	296	C21H44	000629-94-7	90
5	Hexadecane	226	C16H34	000544-76-3	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

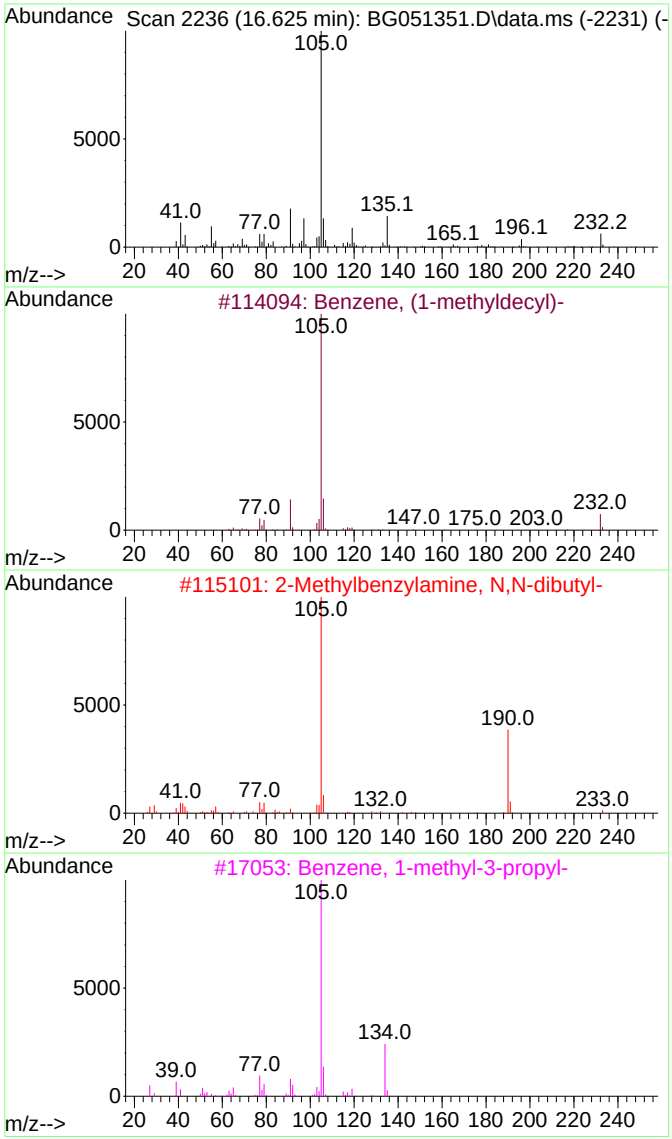
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

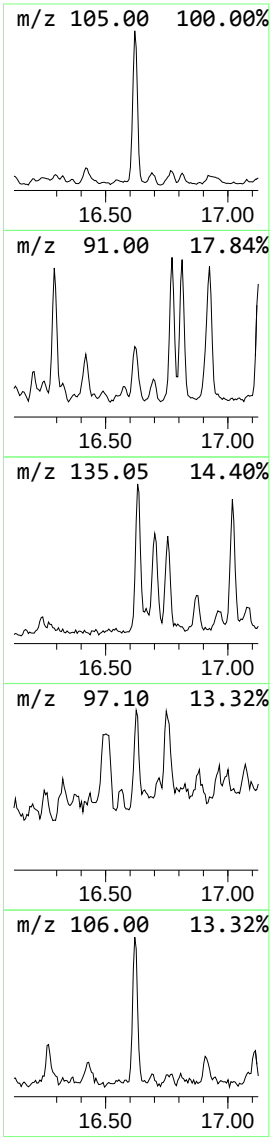
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 Benzene, (1-methyldecyl)- Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.625	17.60 ng/ul	348821	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyldecyl)-	232	C17H28	004536-88-3	60
2	2-Methylbenzylamine, N,N-dibutyl-	233	C16H27N	1000310-39-6	46
3	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	43
4	3-Methylbenzylamine, N,N-dibutyl-	233	C16H27N	1000310-40-5	43
5	Benzene, (1-methylhexyl)-	176	C13H20	002132-84-5	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

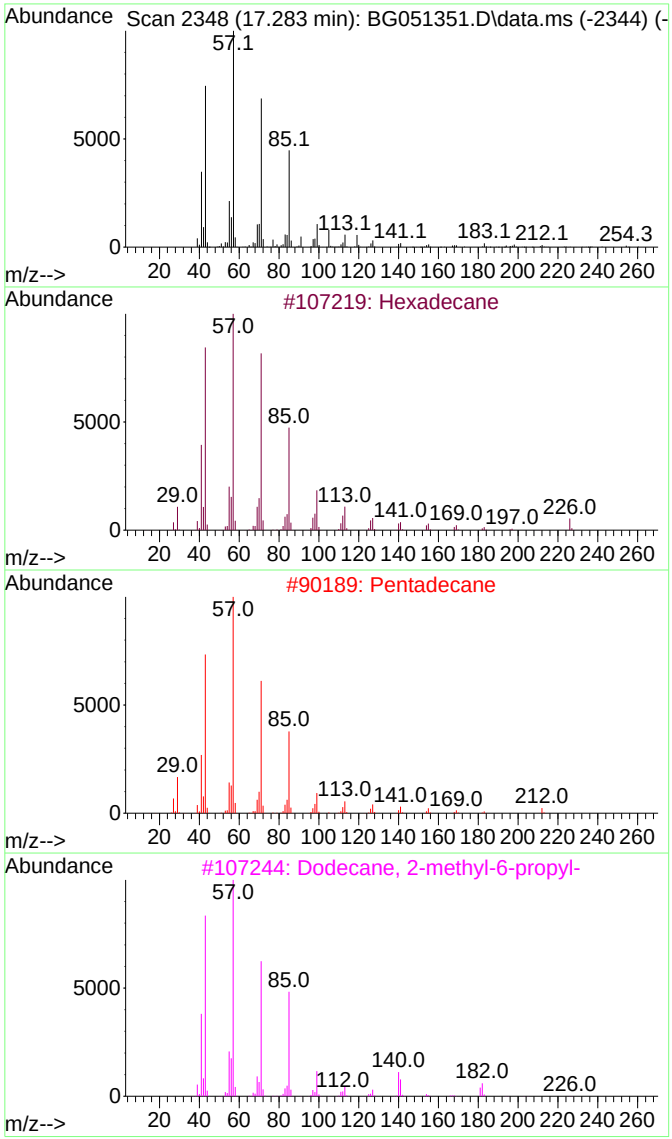
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

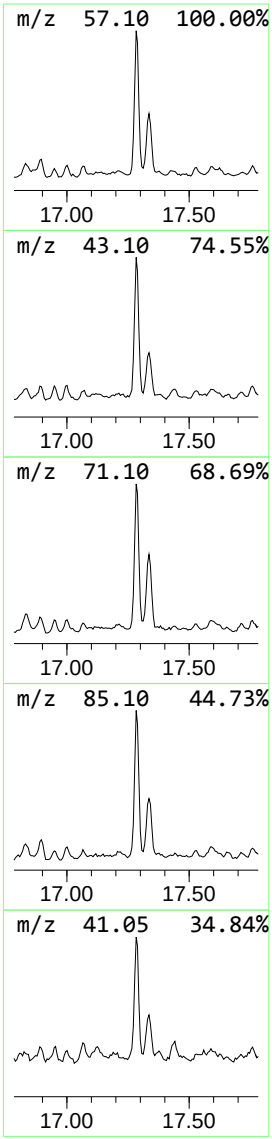
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.283	35.25 ng/ul	698547	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	93
2	Pentadecane	212	C15H32	000629-62-9	93
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4	Tetradecane	198	C14H30	000629-59-4	89
5	Heptadecane	240	C17H36	000629-78-7	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

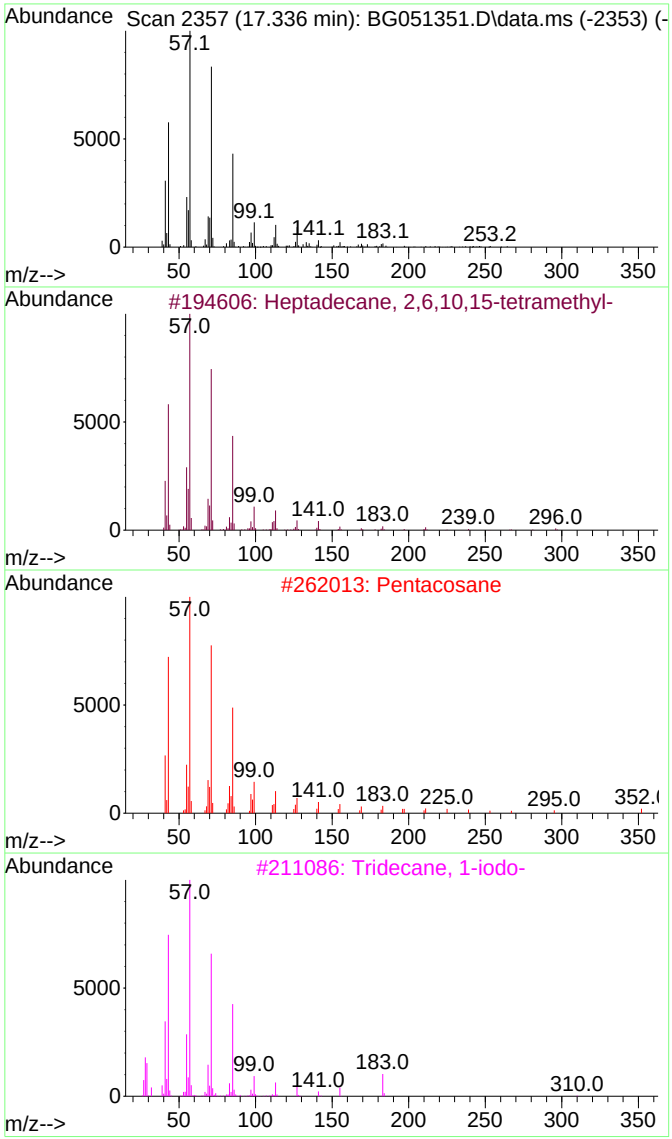
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

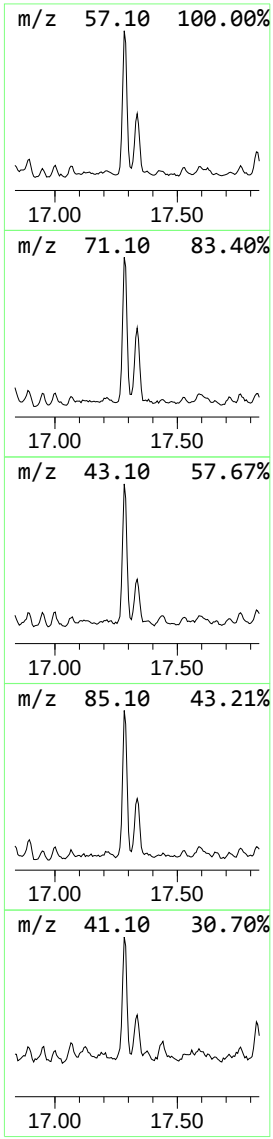
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.336	16.27 ng/ul	322329	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	91
2	Pentacosane	352	C25H52	000629-99-2	90
3	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	90
4	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86
5	Undecane, 2-methyl-	170	C12H26	007045-71-8	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

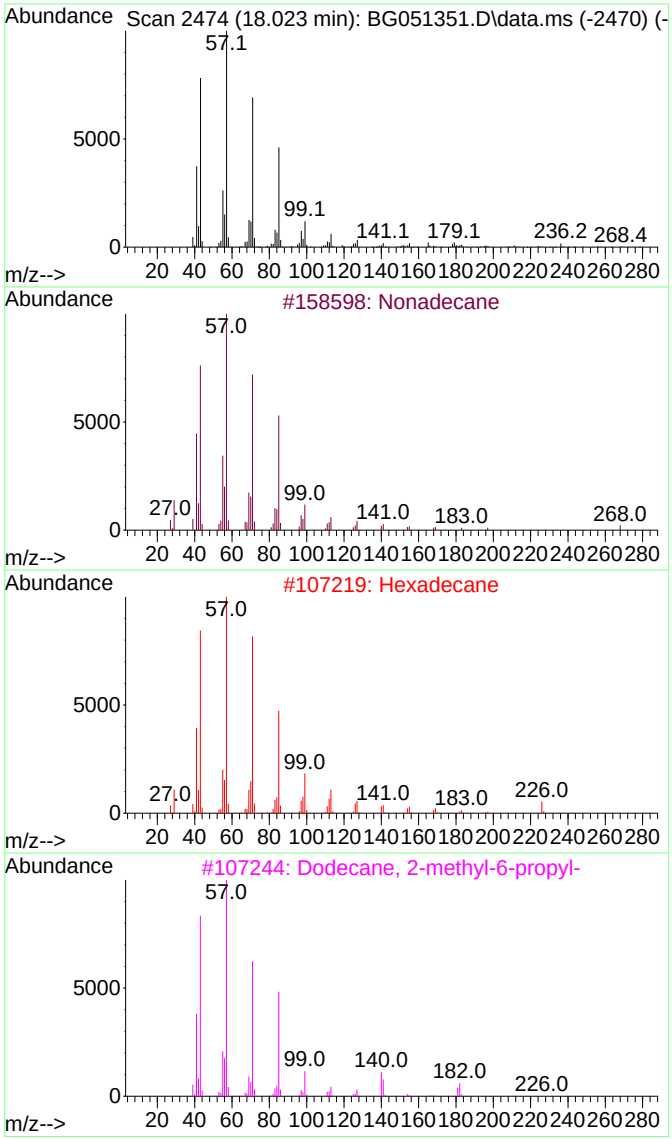
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

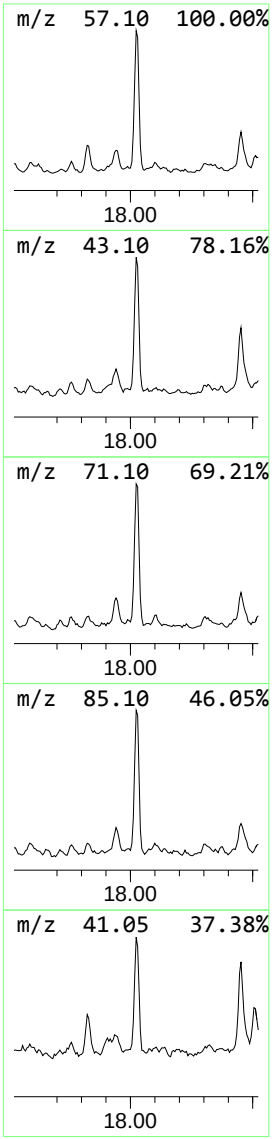
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.023	35.12 ng/ul	695918	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonadecane	268	C19H40	000629-92-5	98
2	Hexadecane	226	C16H34	000544-76-3	94
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
4	Nonadecane	268	C19H40	000629-92-5	91
5	10-Methylnonadecane	282	C20H42	056862-62-5	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

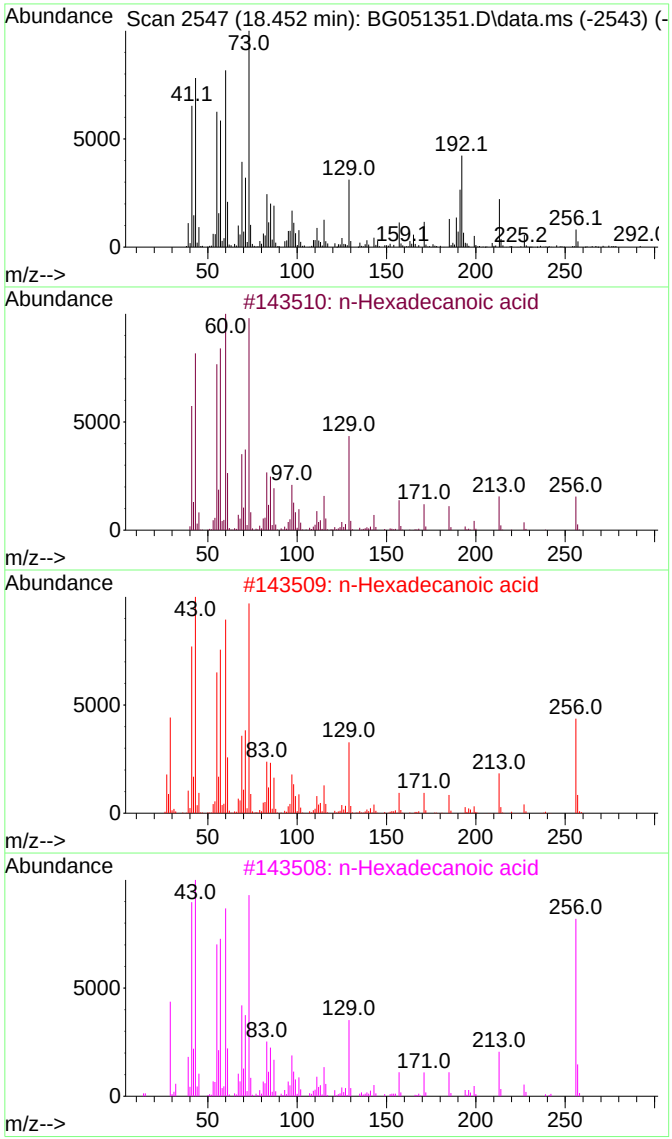
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

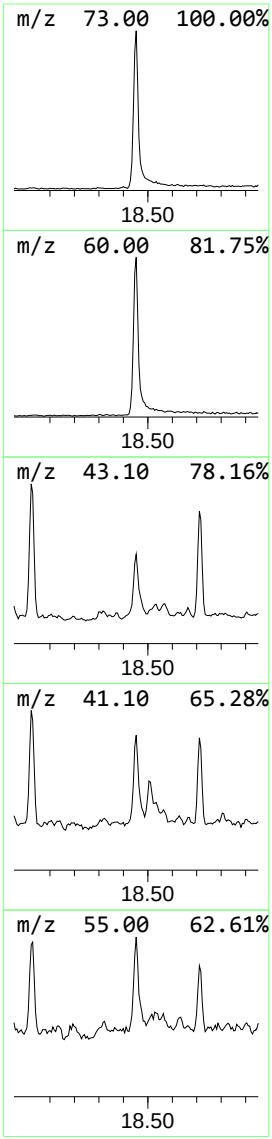
****

Peak Number 48 n-Hexadecanoic acid Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.452	34.13 ng/ul	676355	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	Tridecanoic acid	214	C13H26O2	000638-53-9	91
5	Tridecanoic acid	214	C13H26O2	000638-53-9	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

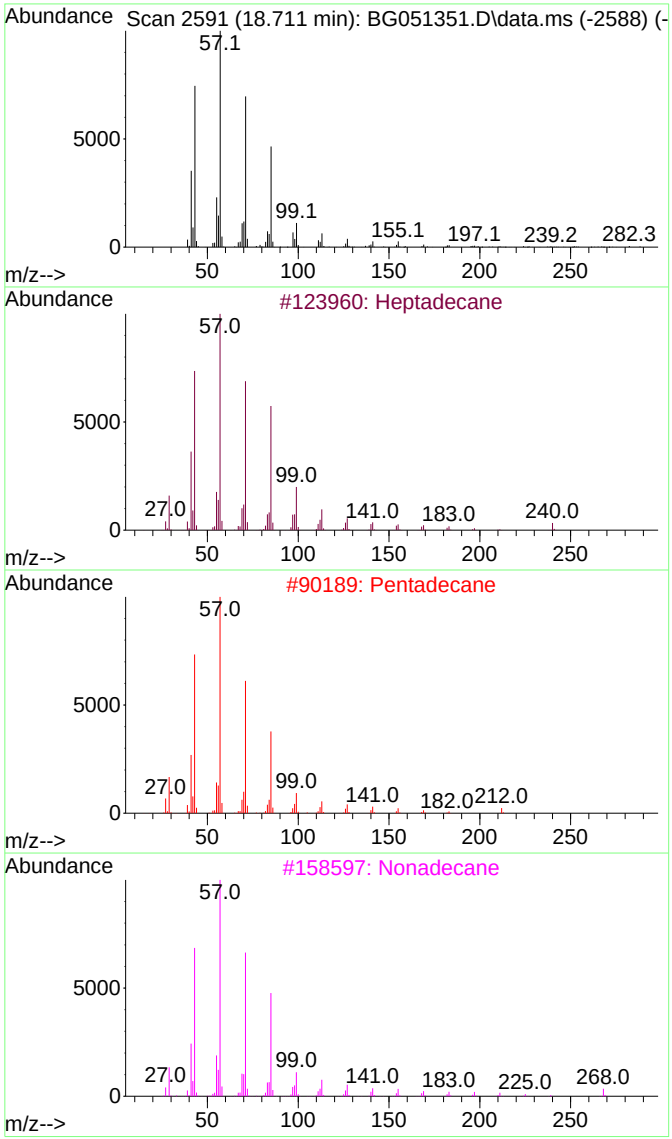
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

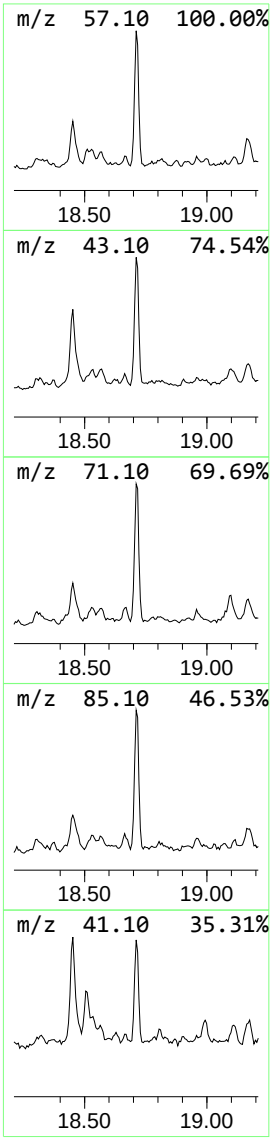
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.711	22.01 ng/ul	436117	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Pentadecane	212	C15H32	000629-62-9	91
3	Nonadecane	268	C19H40	000629-92-5	91
4	Heneicosane	296	C21H44	000629-94-7	90
5	Tetracosane	338	C24H50	000646-31-1	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

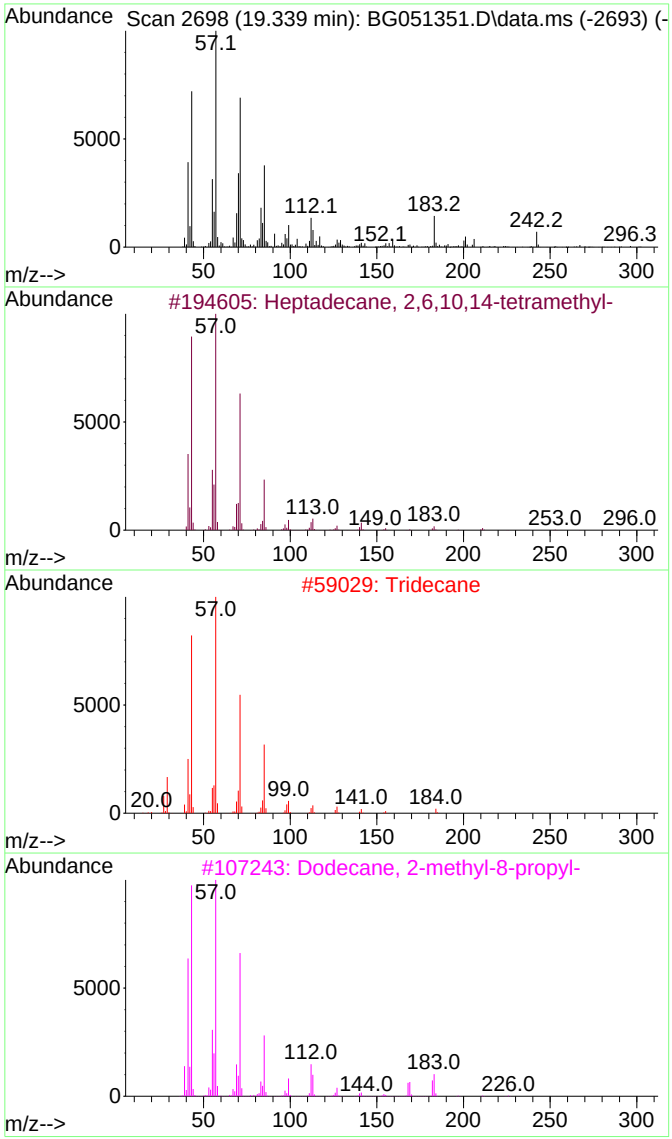
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

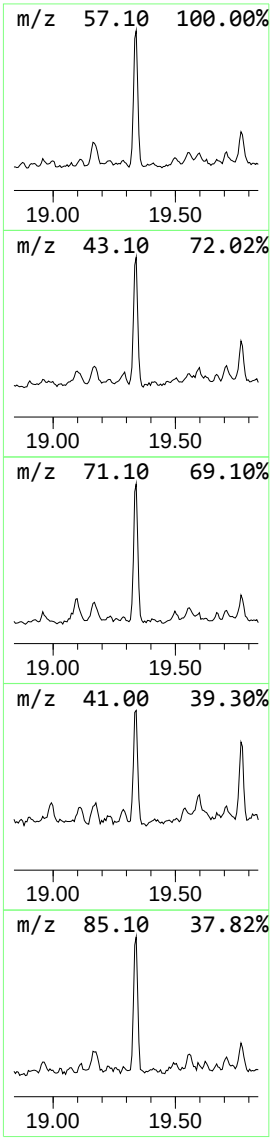
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.339	39.98 ng/ul	792230	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	83
2	Tridecane	184	C13H28	000629-50-5	76
3	Dodecane, 2-methyl-8-propyl-	226	C16H34	055045-07-3	76
4	Pentadecane, 7-methyl-	226	C16H34	006165-40-8	72
5	Tridecane, 7-propyl-	226	C16H34	055045-09-5	70





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

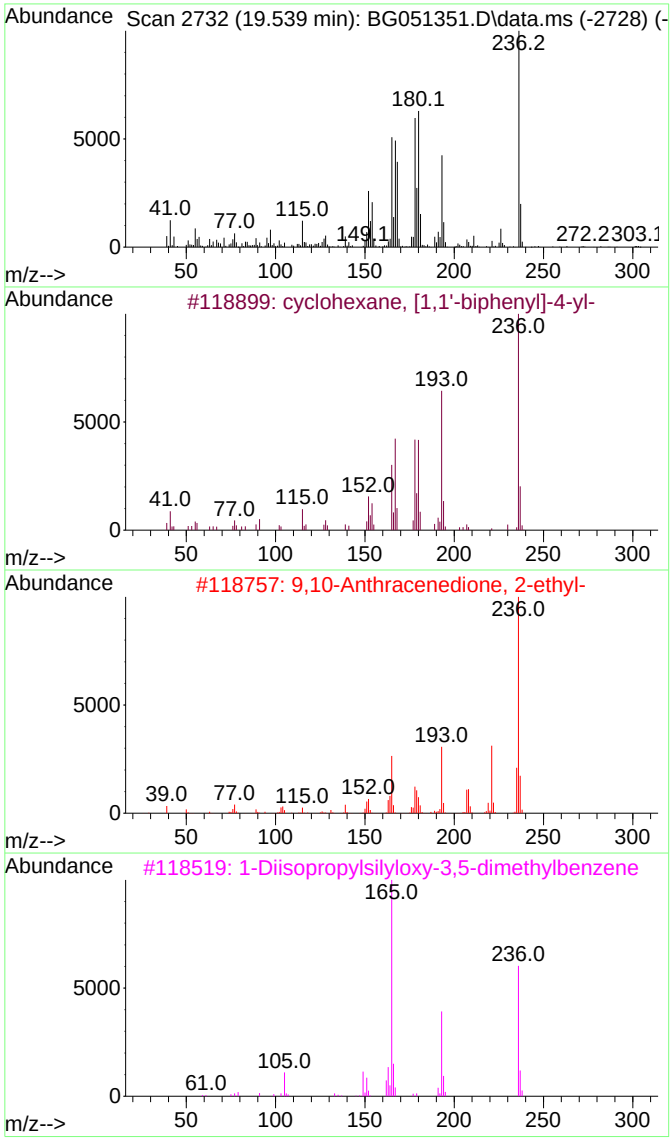
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

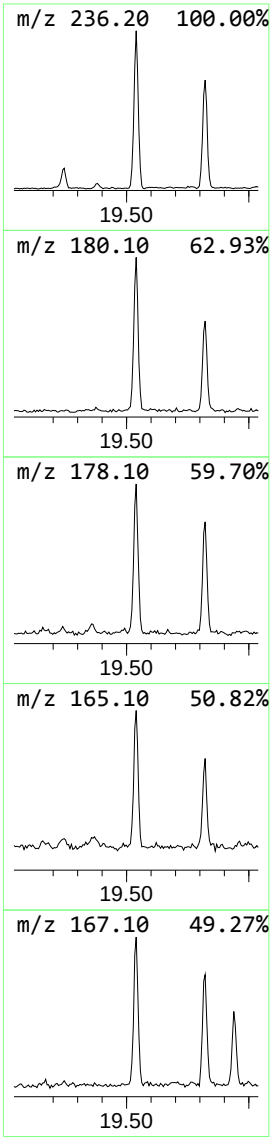
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 52 cyclohexane, [1,1'-biphenyl... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	17.91 ng/ul	354883	Phenanthrene-d10	17.583

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	cyclohexane, [1,1'-biphenyl]-4-yl-	236	C18H20	1000401-12-4	93
2	9,10-Anthracenedione, 2-ethyl-	236	C16H12O2	000084-51-5	46
3	1-Diisopropylsilyloxy-3,5-dimeth...	236	C14H24OSi	1000307-97-3	45
4	9,10-Anthracenedione, 1-ethyl-	236	C16H12O2	024624-29-1	43
5	9,10-Anthracenedione, 2,3-dimethyl-	236	C16H12O2	006531-35-7	38





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

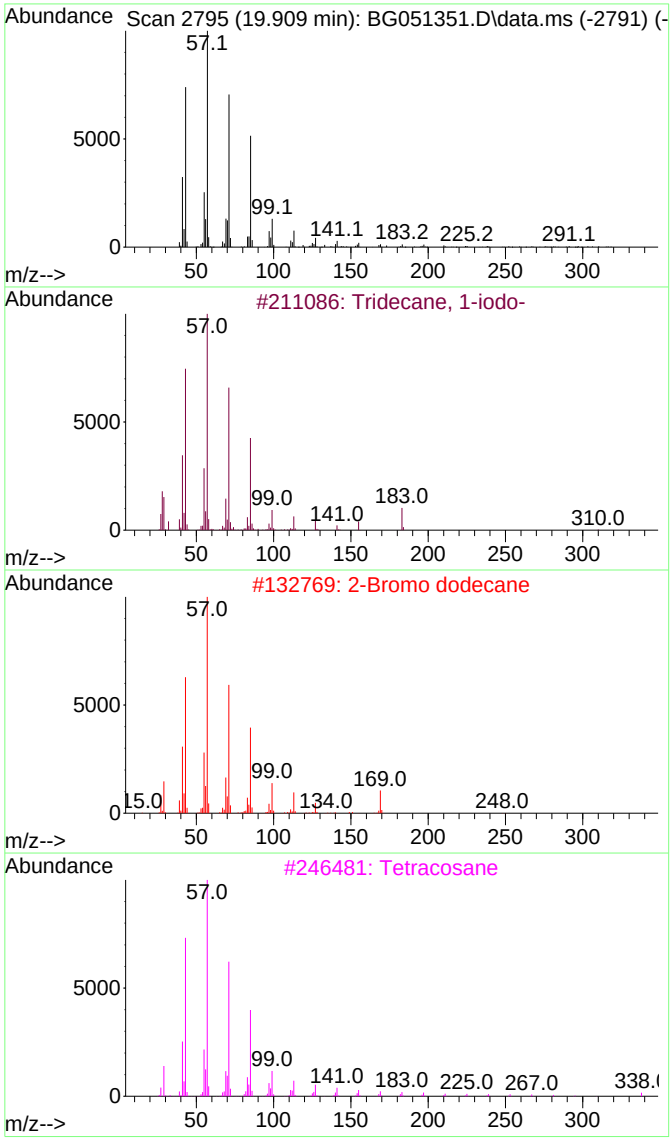
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

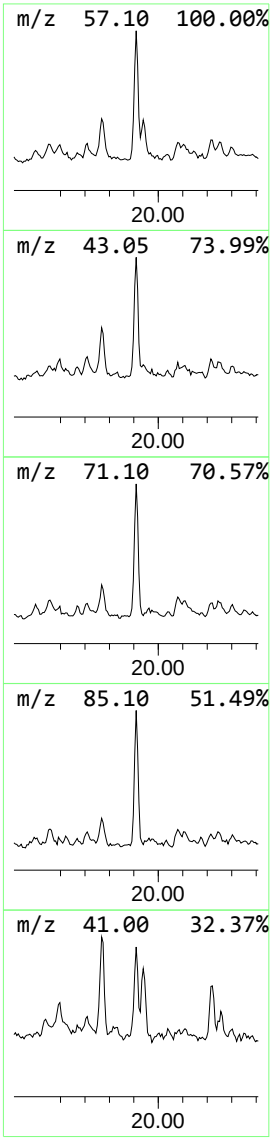
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 55 Tridecane, 1-iodo- Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.909	17.29 ng/ul	372884	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	93
2	2-Bromo dodecane	248	C12H25Br	013187-99-0	90
3	Tetracosane	338	C24H50	000646-31-1	90
4	Tetradecane	198	C14H30	000629-59-4	86
5	Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	86





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

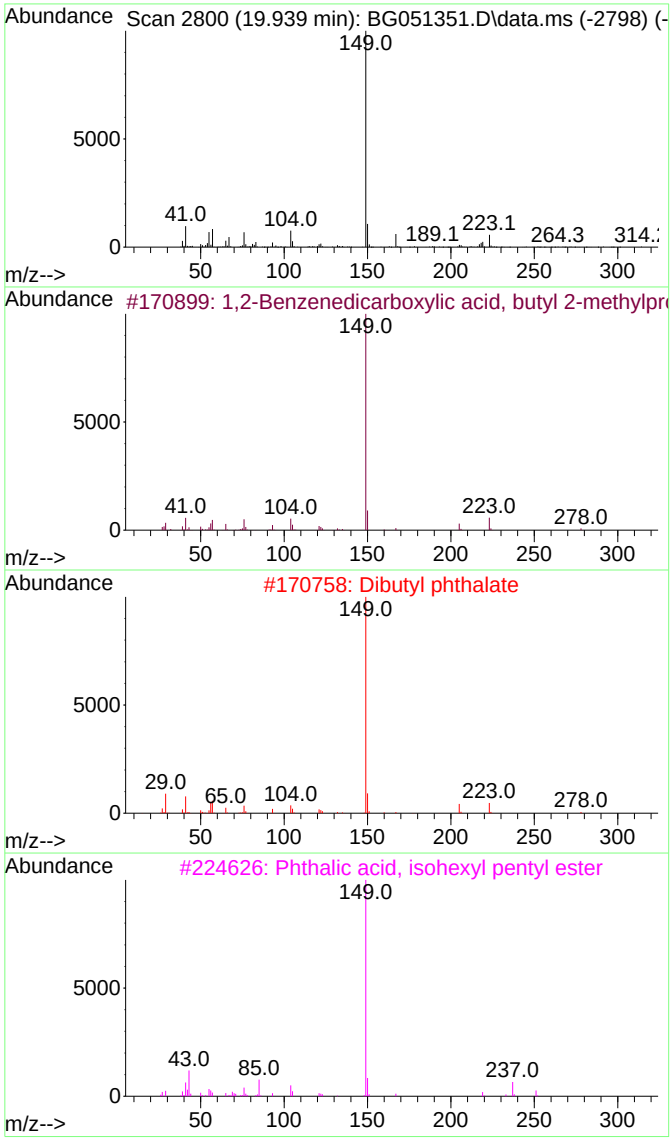
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

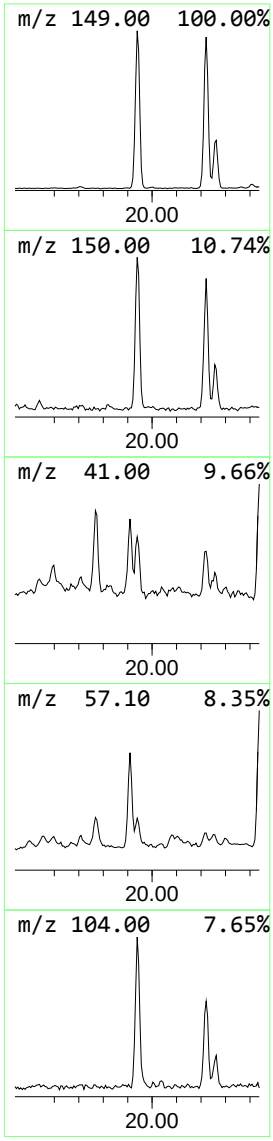
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 56 1,2-Benzenedicarboxylic aci... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.939	23.78 ng/ul	512652	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, bu...	278	C16H22O4	017851-53-5	86
2	Dibutyl phthalate	278	C16H22O4	000084-74-2	86
3	Phthalic acid, isohexyl pentyl e...	320	C19H28O4	1000308-95-8	78
4	Phthalic acid, 8-chlorooctyl iso...	368	C20H29ClO4	1000356-85-8	78
5	Phthalic acid, isobutyl undecyl ...	376	C23H36O4	1010308-97-3	78





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

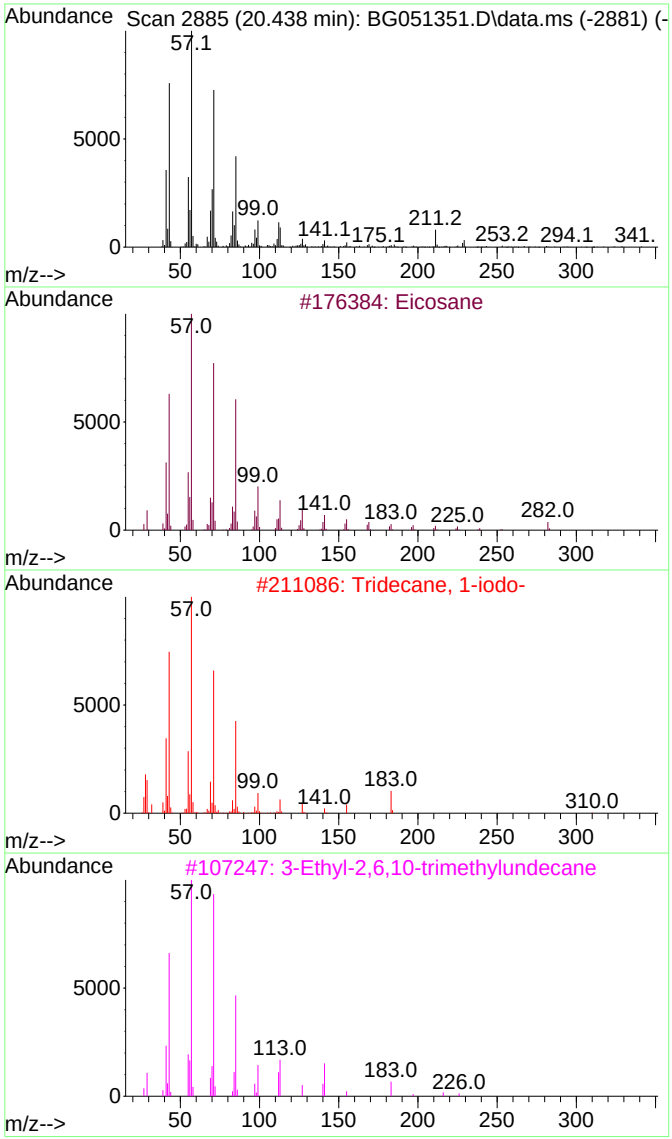
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

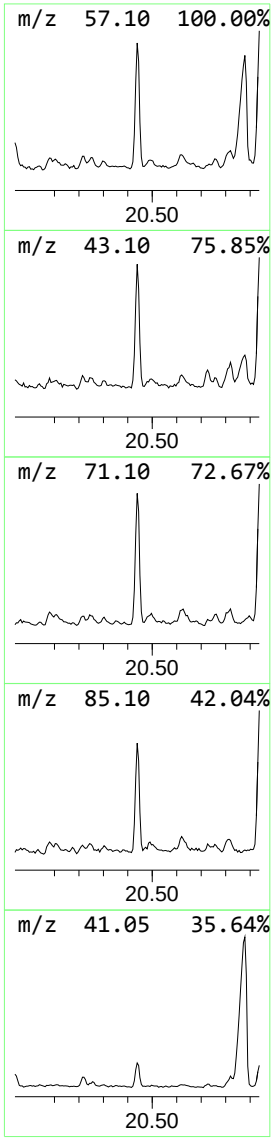
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.438	31.59 ng/ul	681023	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	Tridecane, 1-iodo-	310	C13H27I	035599-77-0	89
3	3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	83
4	Tetradecane	198	C14H30	000629-59-4	80
5	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	74





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

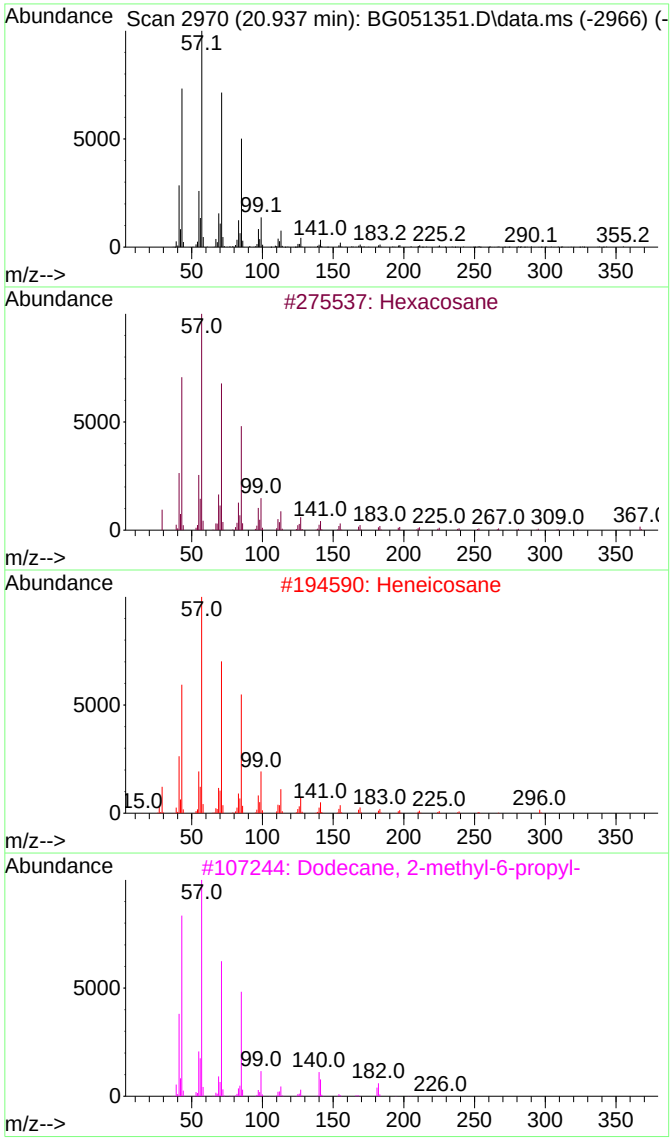
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

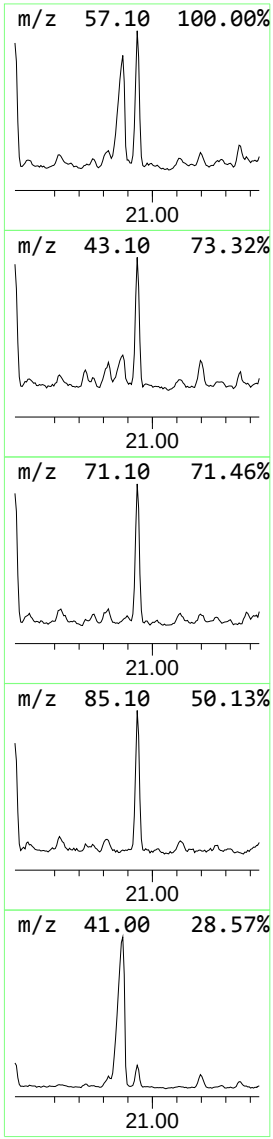
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.937	27.42 ng/ul	591229	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	91
2	Heneicosane	296	C21H44	000629-94-7	91
3	Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	90
4	Octacosane	394	C28H58	000630-02-4	90
5	Nonadecane	268	C19H40	000629-92-5	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

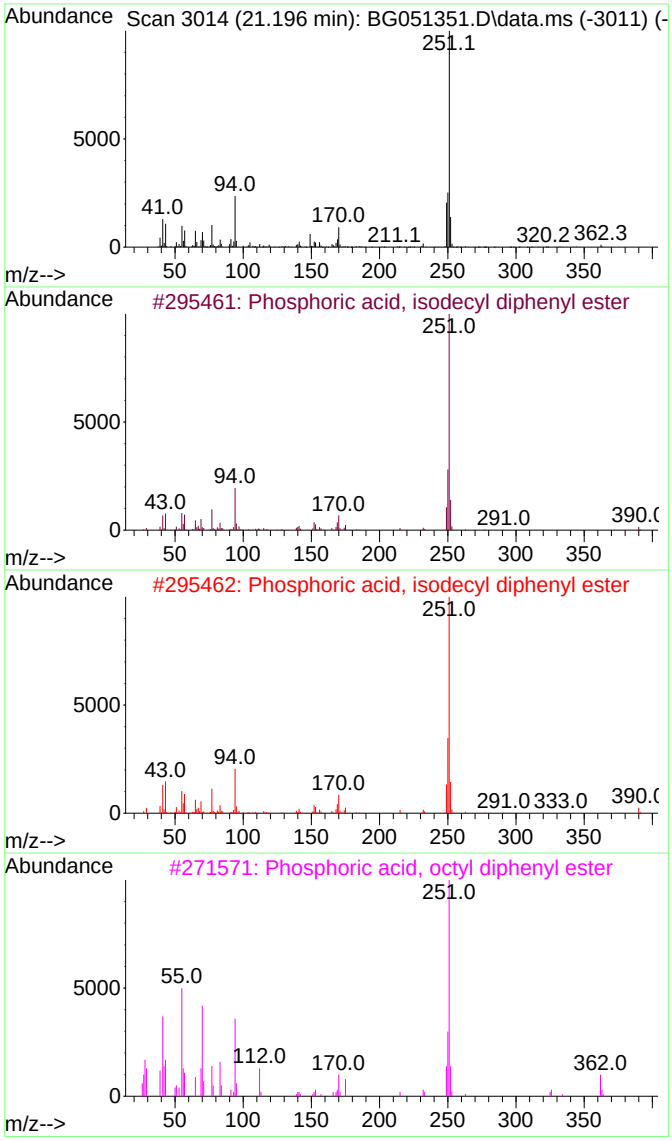
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

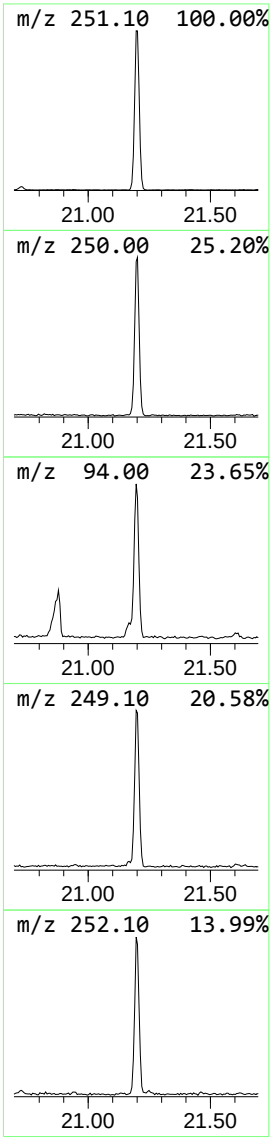
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 60 Phosphoric acid, isodecyl d... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.196	32.63 ng/ul	703489	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phosphoric acid, isodecyl diphen...	390	C22H31O4P	1346599-15-2	87
2	Phosphoric acid, isodecyl diphen...	390	C22H31O4P	1346599-15-2	83
3	Phosphoric acid, octyl diphenyl ...	362	C20H27O4P	000115-88-8	50
4	Octicizer	362	C20H27O4P	001241-94-7	49
5	trans-4-(Dimethylamino)chalcone	251	C17H17NO	022965-98-6	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

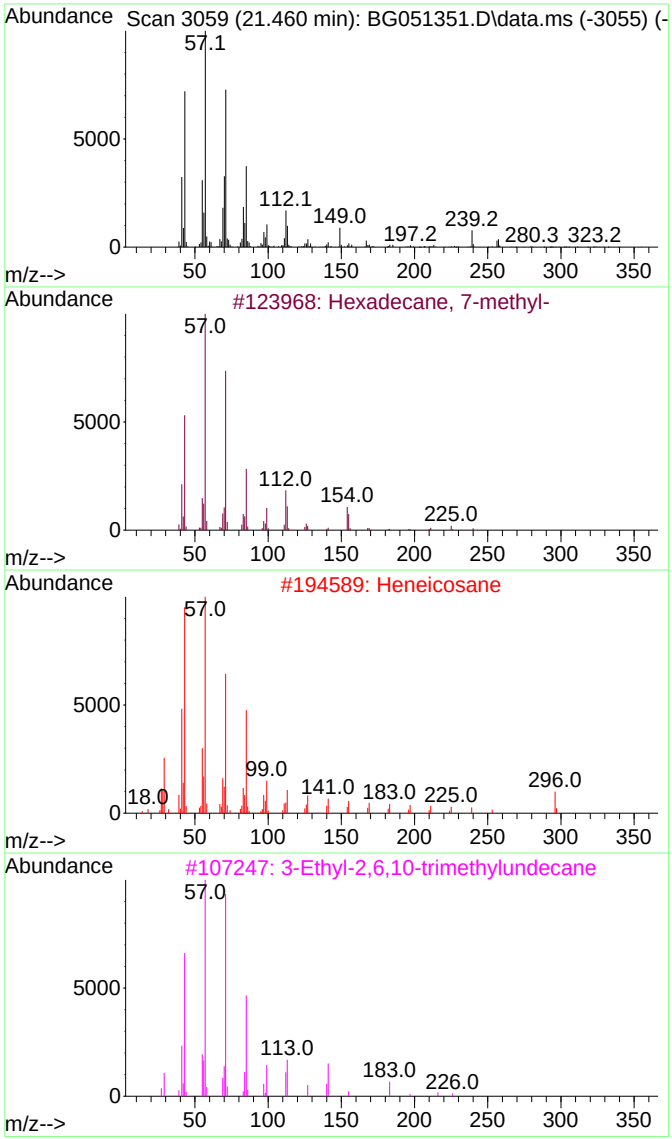
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

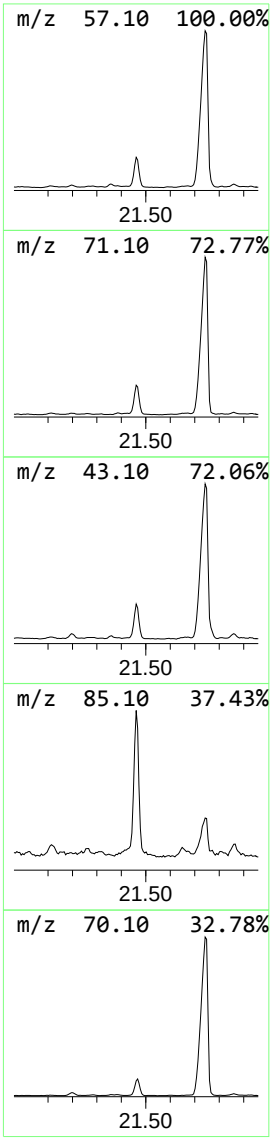
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.460	61.99 ng/ul	1336590	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 7-methyl-	240	C17H36	026730-20-1	94
2	Heneicosane	296	C21H44	000629-94-7	93
3	3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	87
4	Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	86
5	Dodecane	170	C12H26	000112-40-3	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

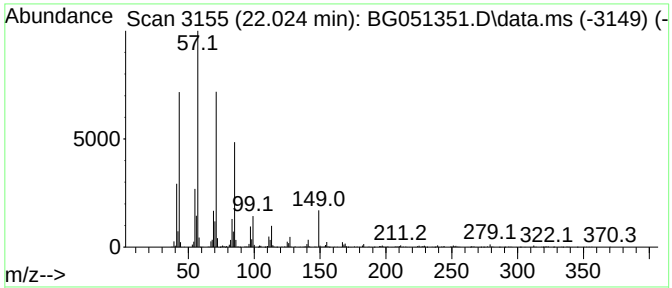
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

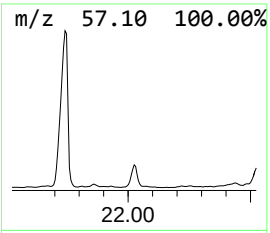
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

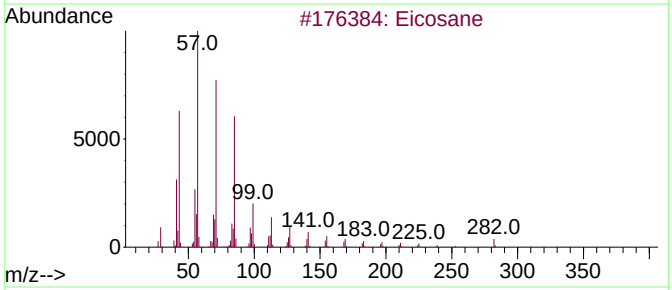
****
Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 15

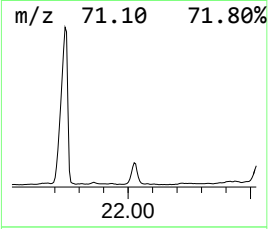
R.T.	EstConc	Area	Relative to ISTD	R.T.
22.024	52.50 ng/ul	1132070	Chrysene-d12	21.883

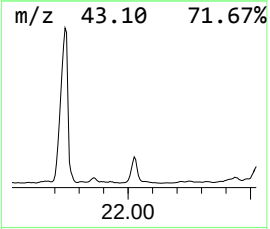
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Eicosane	282	C20H42	000112-95-8	95
2	Octacosane	394	C28H58	000630-02-4	87
3	Heneicosane	296	C21H44	000629-94-7	87
4	Hexacosane	366	C26H54	000630-01-3	87
5	Tetracosane	338	C24H50	000646-31-1	87

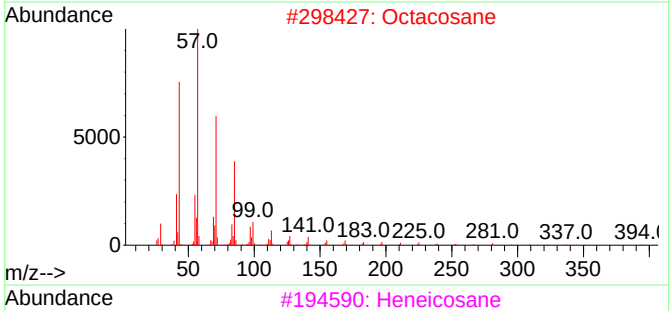


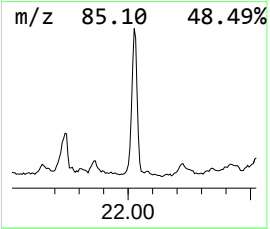


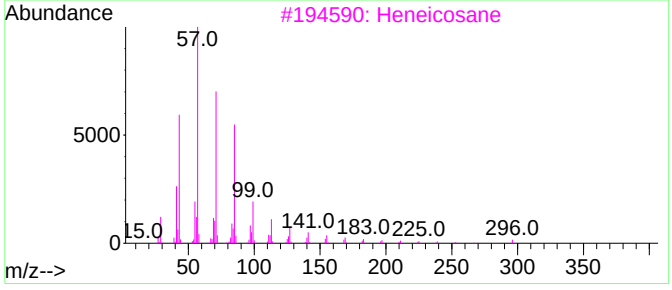


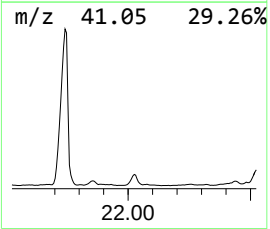












Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

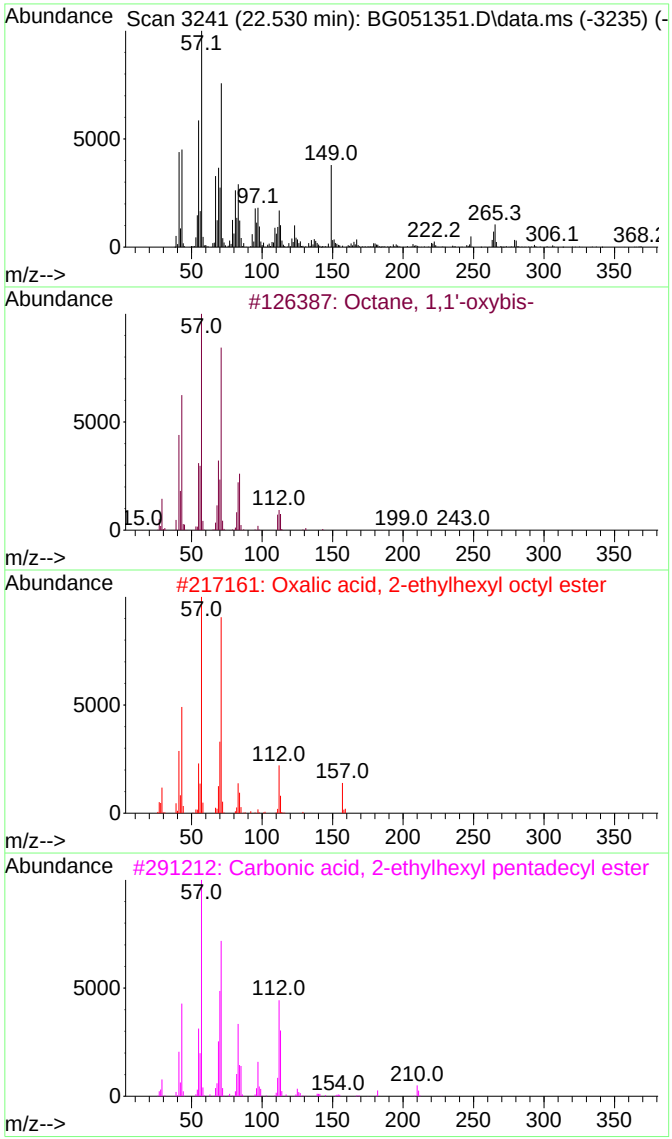
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

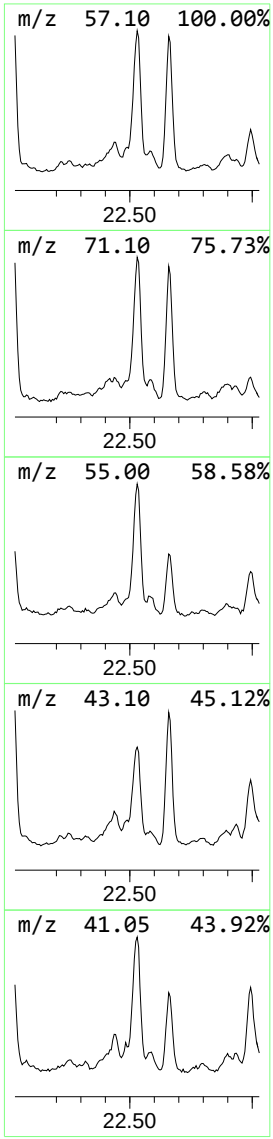
****

Peak Number 65 Octane, 1,1'-oxybis- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.530	100.04 ng/ul	2157010	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	52
2	Oxalic acid, 2-ethylhexyl octyl ...	314	C18H34O4	1000309-39-1	47
3	Carbonic acid, 2-ethylhexyl pent...	384	C24H48O3	1000383-14-2	46
4	Hexadecane, 7-methyl-	240	C17H36	026730-20-1	43
5	Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	43





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

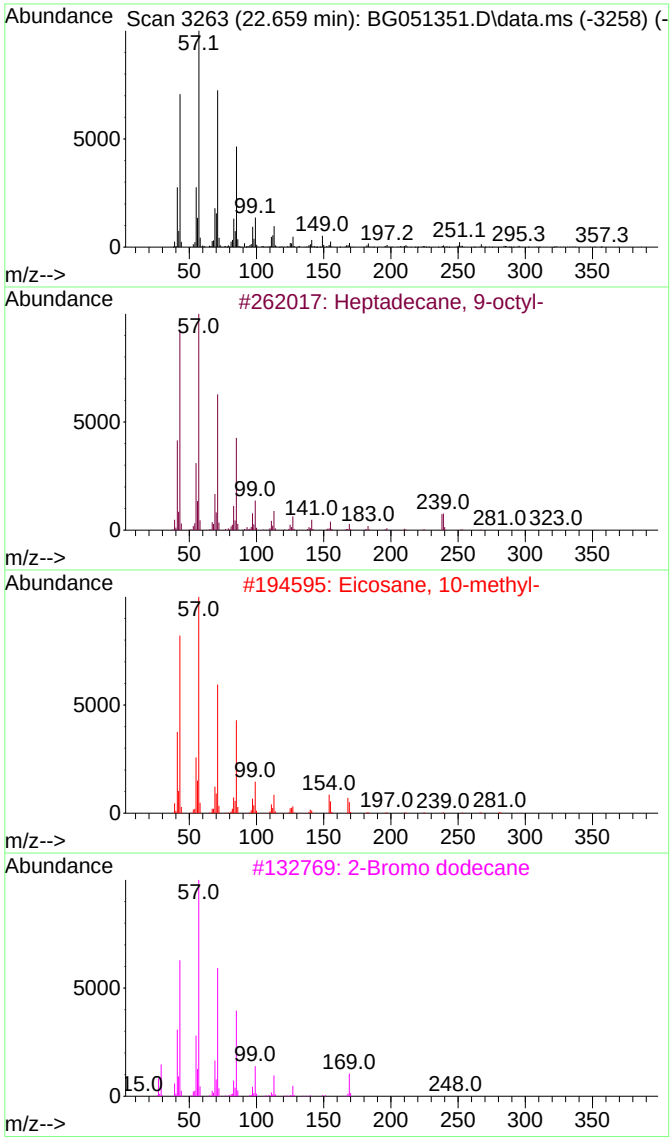
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

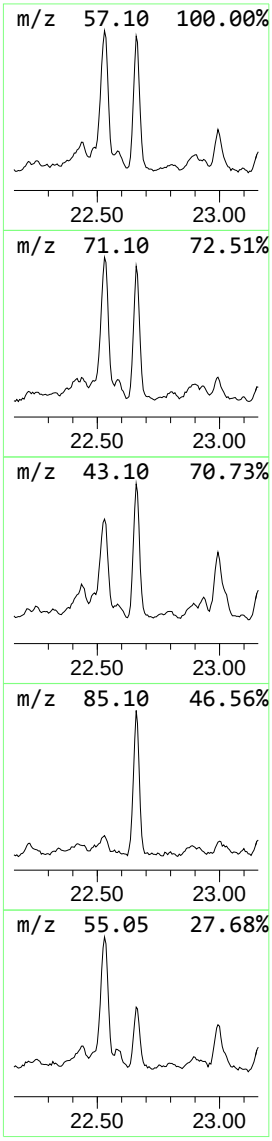
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 66 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.659	52.80 ng/ul	1138360	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
2	Eicosane, 10-methyl-	296	C21H44	054833-23-7	91
3	2-Bromo dodecane	248	C12H25Br	013187-99-0	91
4	Hexadecane	226	C16H34	000544-76-3	91
5	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

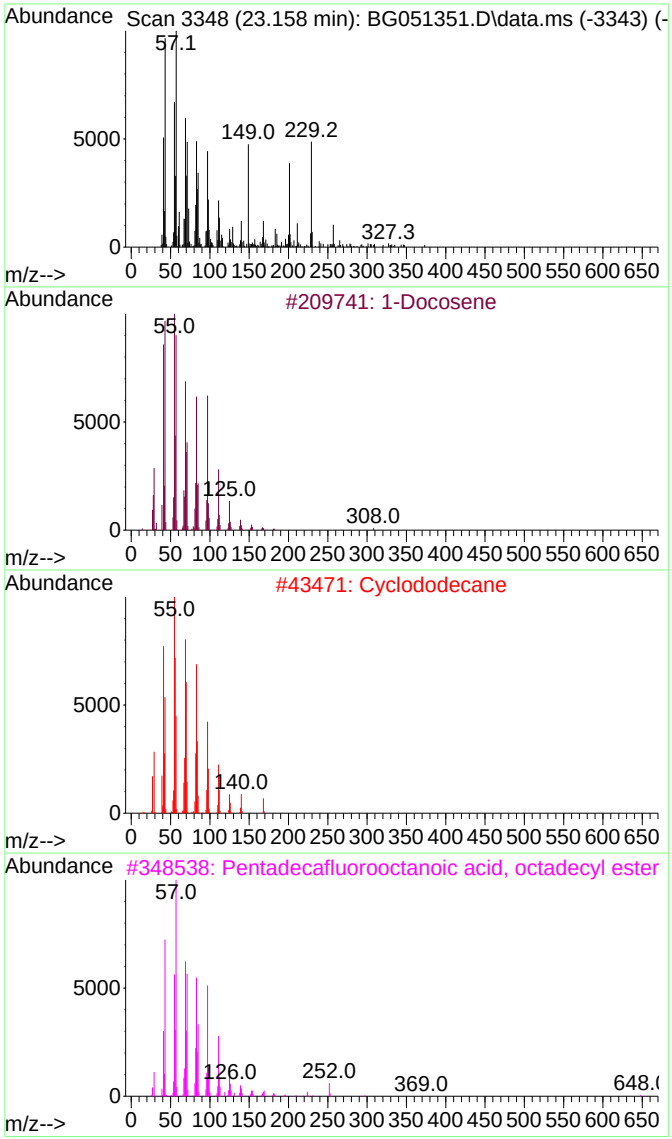
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

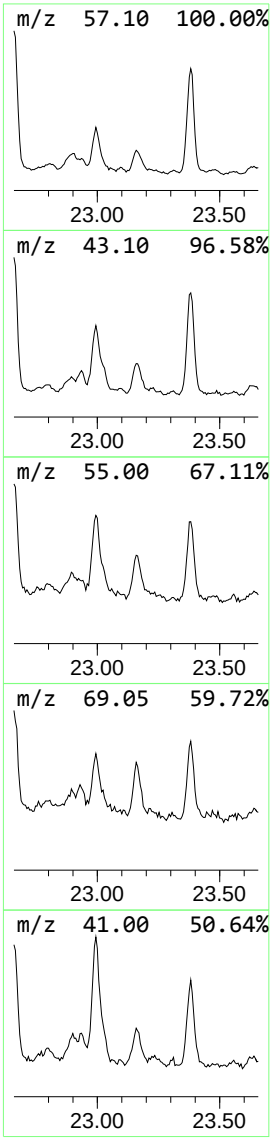
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.158	20.27 ng/ul	437064	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	86
2	Cyclododecane	168	C12H24	000294-62-2	74
3	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	55
4	3-Eicosene, (E)-	280	C20H40	074685-33-9	50
5	Dichloroacetic acid, tetradecyl ...	324	C16H30Cl2O2	083005-02-1	50





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

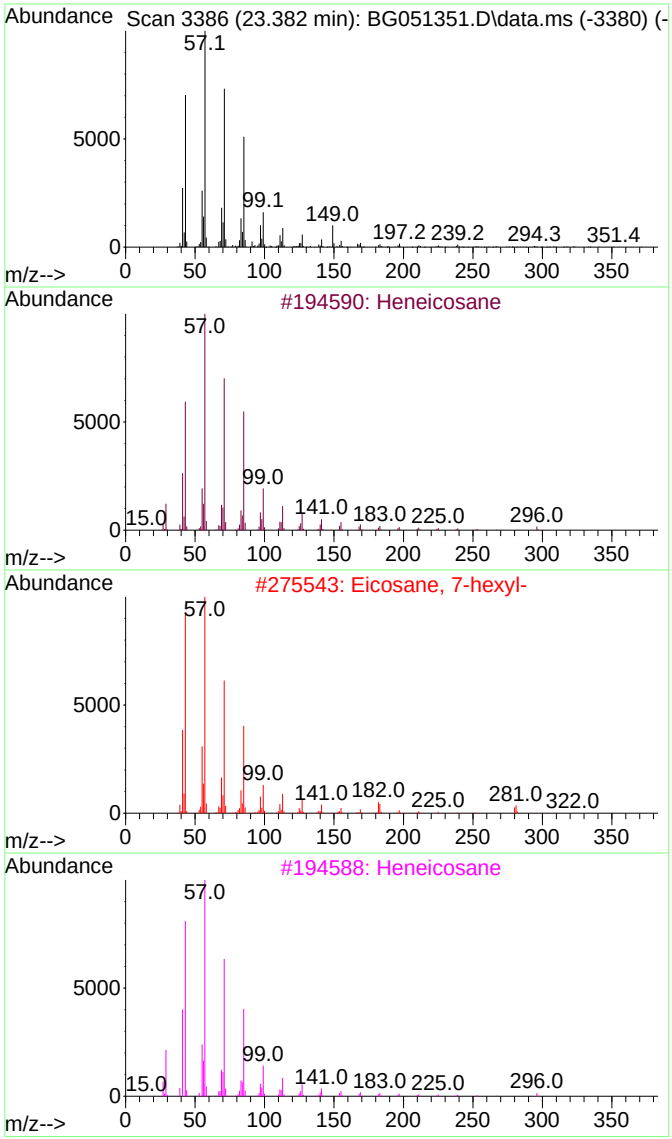
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

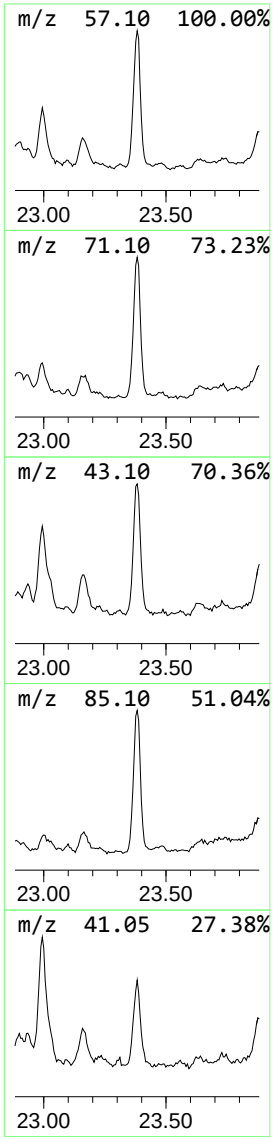
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.382	44.44 ng/ul	958119	Chrysene-d12	21.883

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heneicosane	296	C21H44	000629-94-7	91
2	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	91
3	Heneicosane	296	C21H44	000629-94-7	90
4	Octacosane	394	C28H58	000630-02-4	87
5	Heptacosane	380	C27H56	000593-49-7	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

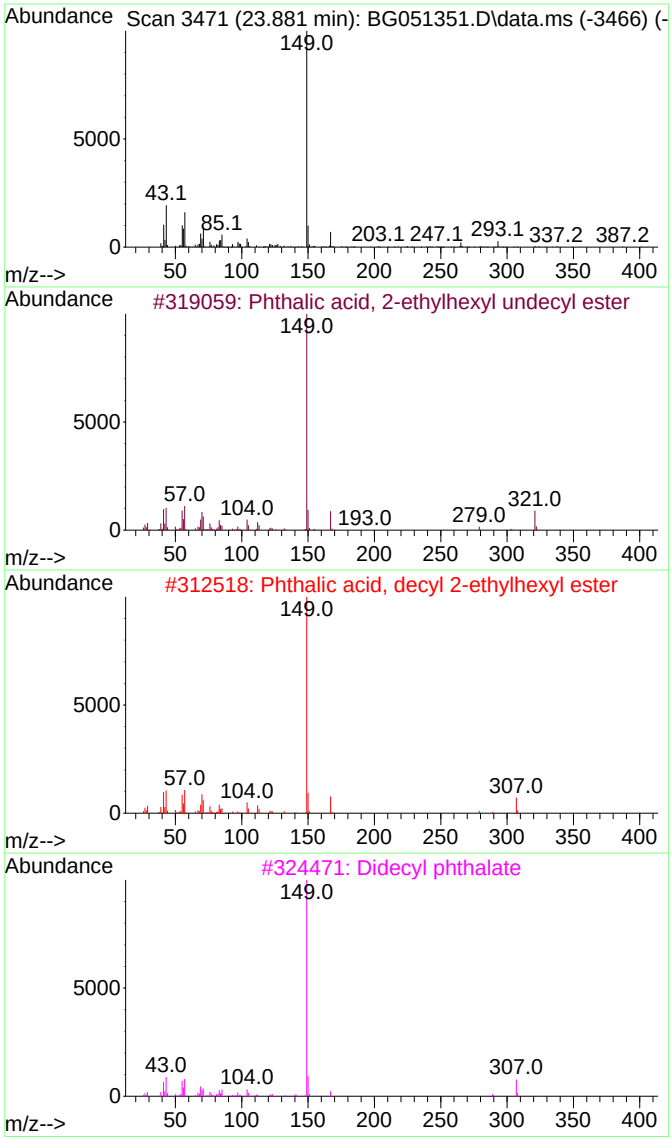
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

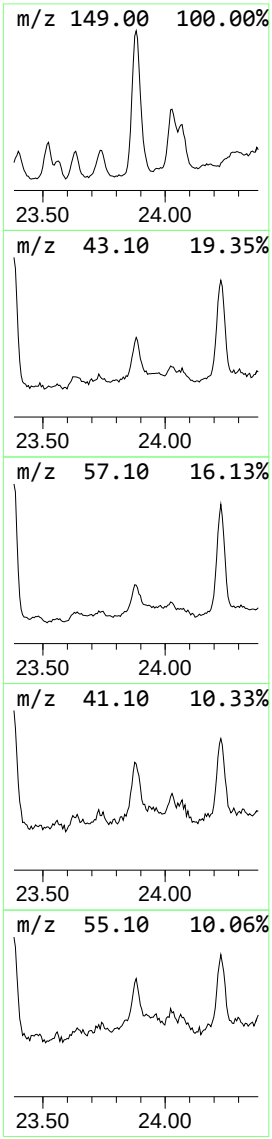
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 69 Phthalic acid, 2-ethylhexyl... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.881	20.88 ng/ul	472290	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phthalic acid, 2-ethylhexyl unde...	432	C27H44O4	1000308-97-4	86
2	Phthalic acid, decyl 2-ethylhexy...	418	C26H42O4	1000309-00-0	86
3	Didecyl phthalate	446	C28H46O4	000084-77-5	86
4	1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	80
5	Phthalic acid, 3-fluorophenyl he...	498	C31H43FO4	1000356-66-2	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

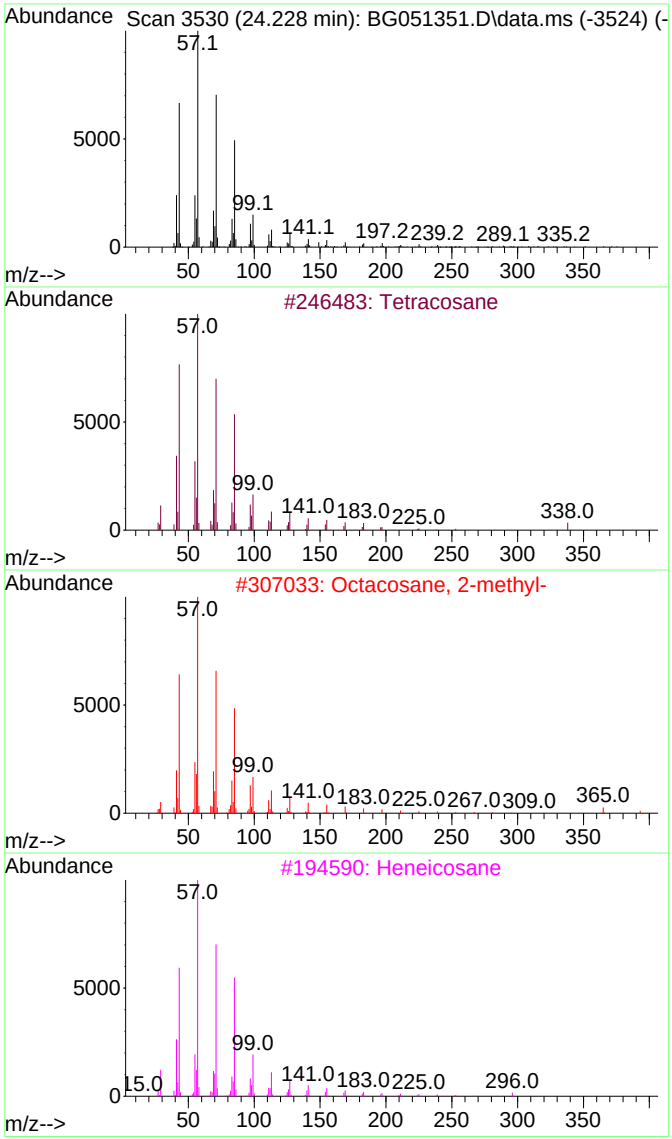
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

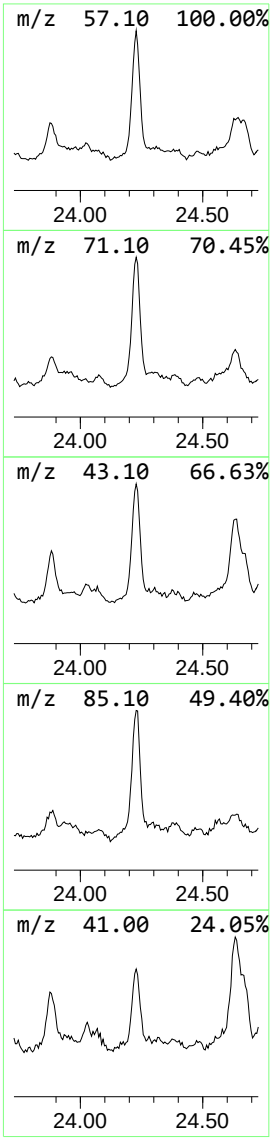
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.228	28.79 ng/ul	651197	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
3	Heneicosane	296	C21H44	000629-94-7	91
4	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	91
5	Hexacosane	366	C26H54	000630-01-3	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

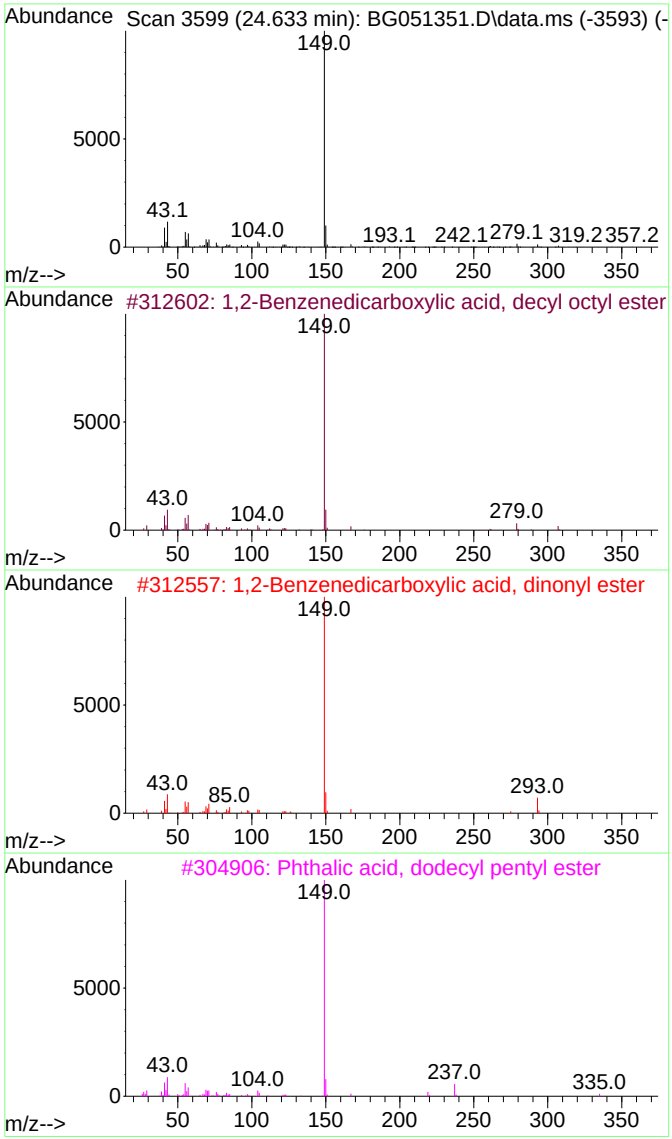
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

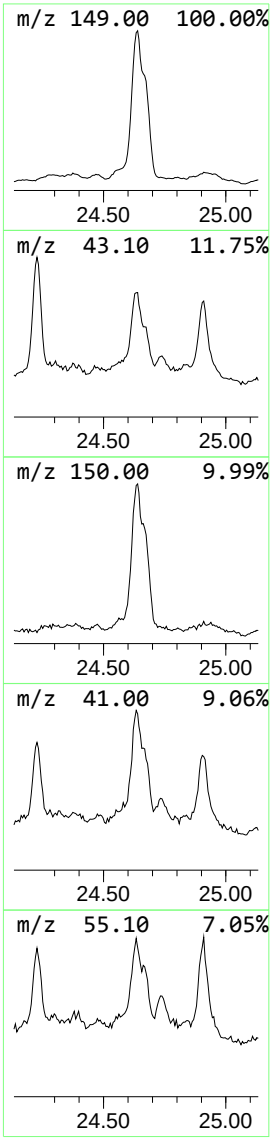
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 71 1,2-Benzenedicarboxylic aci... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.633	71.12 ng/ul	1608510	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	91
2	1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	80
3	Phthalic acid, dodecyl pentyl ester	404	C25H40O4	1000308-93-0	80
4	Phthalic acid, decyl hex-3-yl ester	390	C24H38O4	1000356-96-1	80
5	Di-n-octyl phthalate	390	C24H38O4	000117-84-0	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

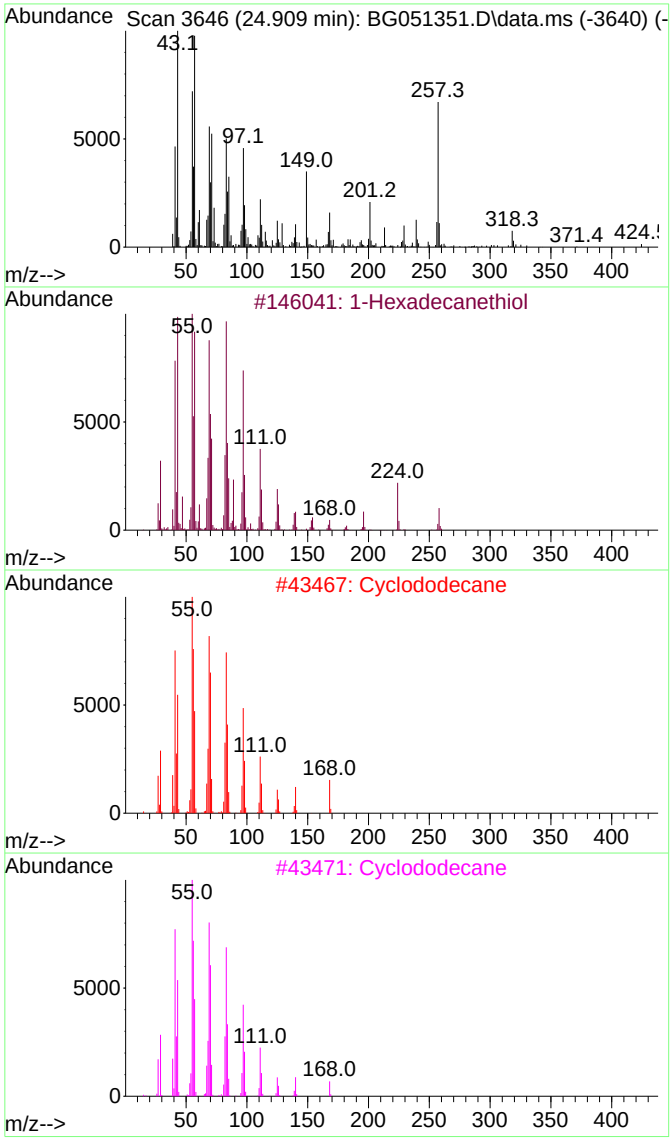
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

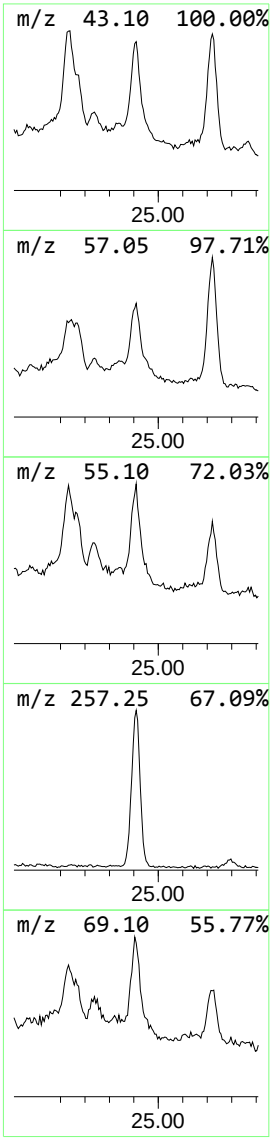
****

Peak Number 72 1-Hexadecanethiol Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.909	45.96 ng/ul	1039500	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexadecanethiol	258	C16H34S	002917-26-2	91
2	Cyclododecane	168	C12H24	000294-62-2	87
3	Cyclododecane	168	C12H24	000294-62-2	86
4	Hexadecanoic acid, decyl ester	396	C26H52O2	042232-27-9	58
5	Acetic acid, trifluoro-, dodecyl...	282	C14H25F3O2	006222-00-0	55





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

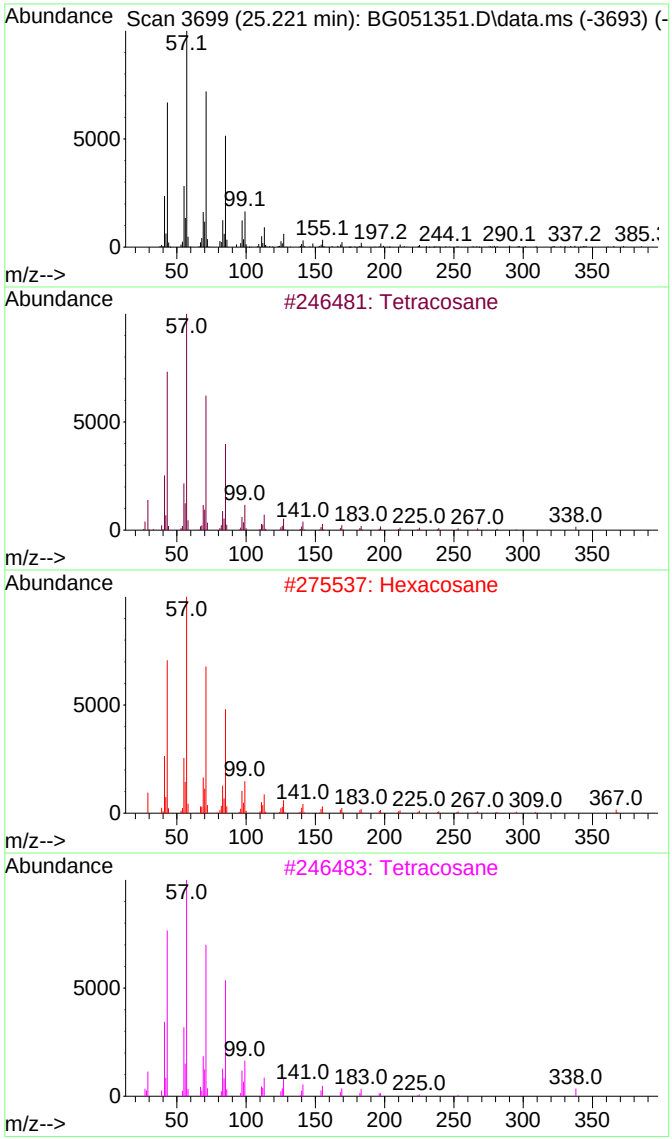
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

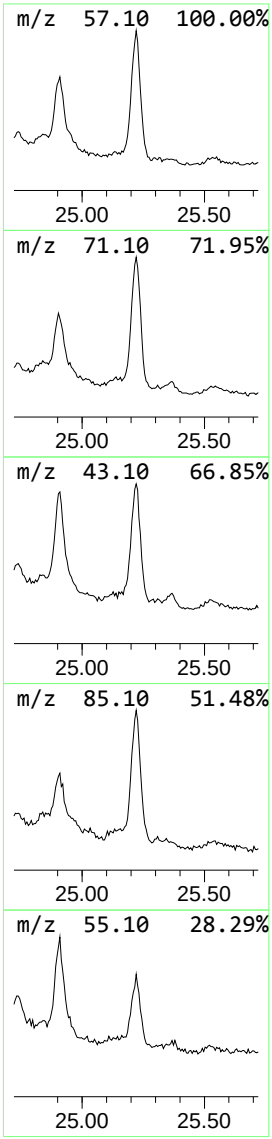
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 73 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.221	20.00 ng/ul	452308	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	91
2	Hexacosane	366	C26H54	000630-01-3	91
3	Tetracosane	338	C24H50	000646-31-1	90
4	Eicosane, 7-hexyl-	366	C26H54	055333-99-8	90
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

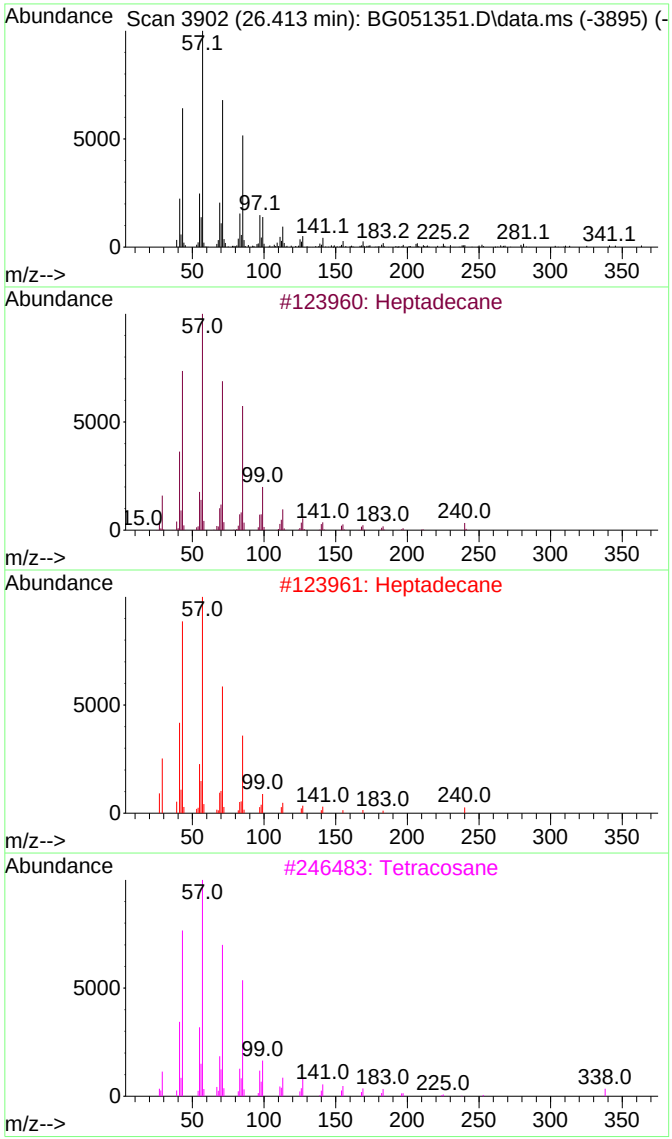
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

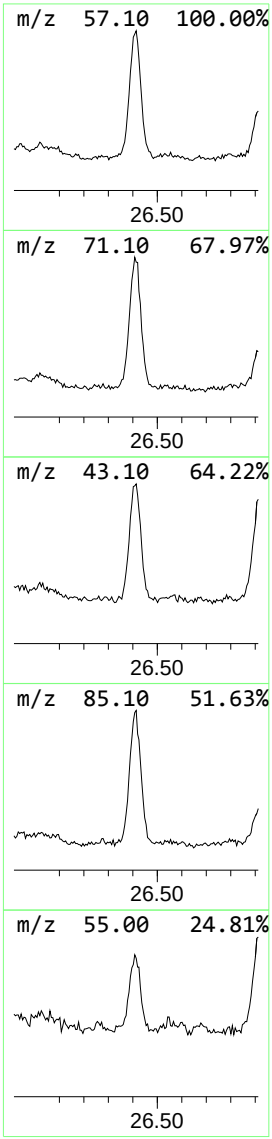
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 74 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.413	24.79 ng/ul	560618	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	95
2	Heptadecane	240	C17H36	000629-78-7	94
3	Tetracosane	338	C24H50	000646-31-1	90
4	Heneicosane	296	C21H44	000629-94-7	87
5	Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	83





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

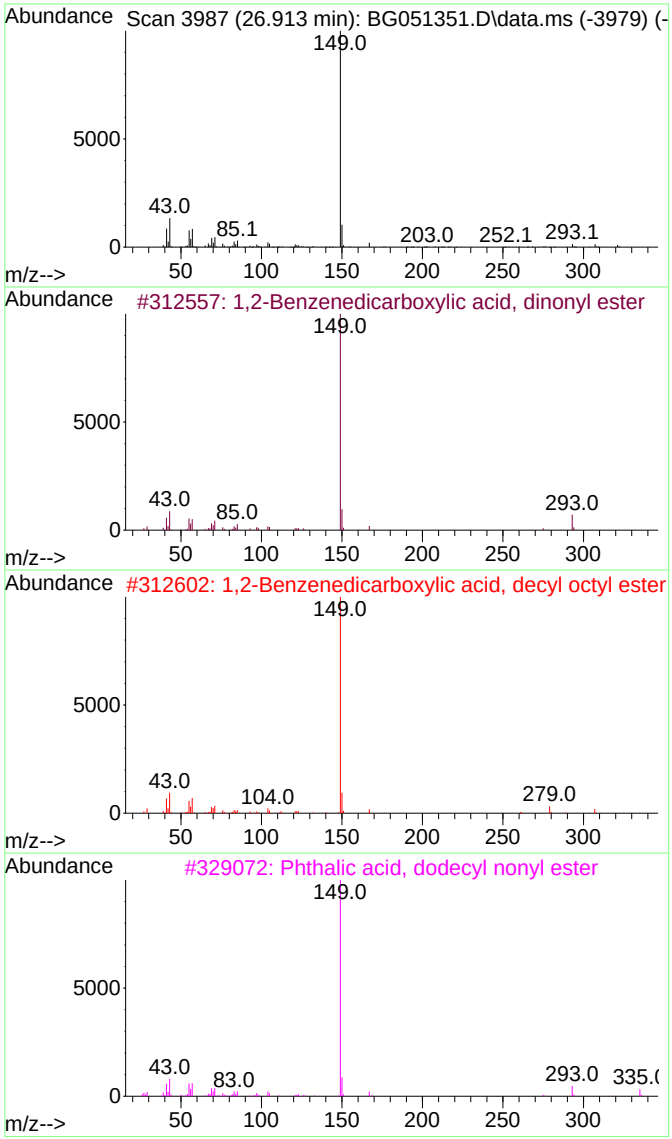
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

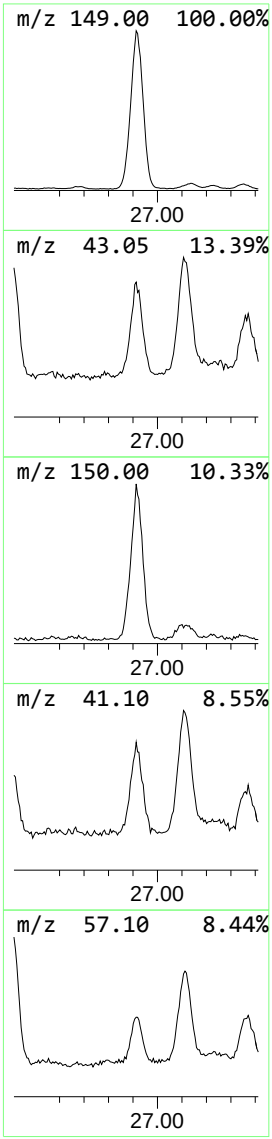
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 75 1,2-Benzenedicarboxylic aci... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.913	42.37 ng/ul	958311	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	83
2	1,2-Benzenedicarboxylic acid, de...	418	C26H42O4	000119-07-3	80
3	Phthalic acid, dodecyl nonyl ester	460	C29H48O4	1000308-92-6	80
4	Phthalic acid, 2-ethylbutyl unde...	404	C25H40O4	1000356-89-6	72
5	Phthalic acid, pentyl pentadecyl...	446	C28H46O4	1000308-92-9	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

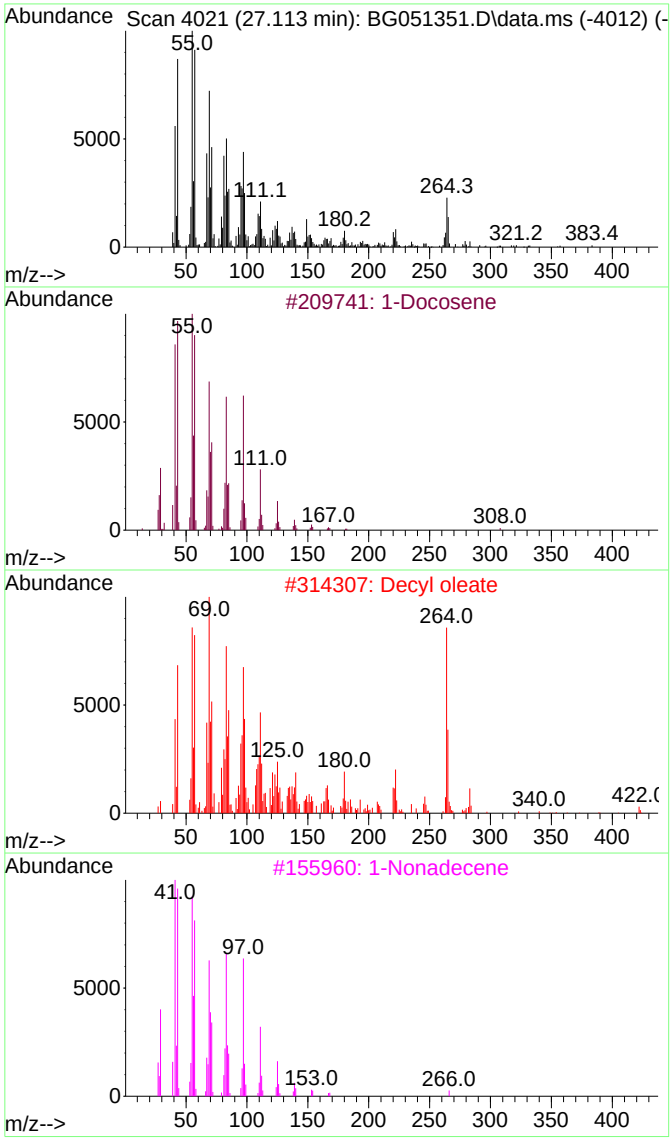
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

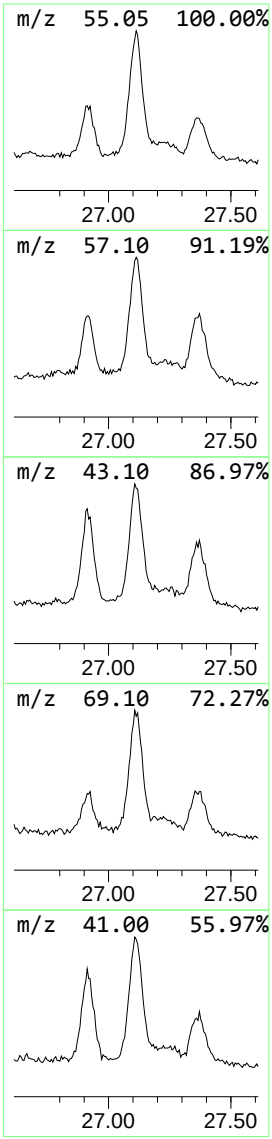
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 76 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.113	53.08 ng/ul	1200350	Perylene-d12	25.291

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	95
2	Decyl oleate	422	C28H54O2	003687-46-5	91
3	1-Nonadecene	266	C19H38	018435-45-5	91
4	9-Octadecenoic acid (Z)-, octade...	535	C36H70O2	017673-49-3	90
5	6-Octadecenoic acid, (Z)-	282	C18H34O2	000593-39-5	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG120621\
Data File : BG051351.D
Acq On : 6 Dec 2021 15:30
Operator : CG/JU
Sample : M4942-03DL 25X
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG112321.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1,3-di...	5.790	113.6	ng/ul	1101820	1	8.194	193936	20.0
Benzene, (1-met...	6.648	16.7	ng/ul	161946	1	8.194	193936	20.0
unknown-01	7.142	30.7	ng/ul	297988	1	8.194	193936	20.0
(DEL) Alkane: S...	7.195	18.3	ng/ul	177414	1	8.194	193936	20.0
Benzene, 1-ethy...	7.253	40.2	ng/ul	389817	1	8.194	193936	20.0
Benzene, 1-ethy...	7.312	42.3	ng/ul	409855	1	8.194	193936	20.0
Benzene, 1-ethy...	7.559	18.9	ng/ul	182869	1	8.194	193936	20.0
(DEL) Alkane: S...	7.776	97.9	ng/ul	949007	1	8.194	193936	20.0
Mesitylene	7.823	77.8	ng/ul	754702	1	8.194	193936	20.0
(DEL) Alkane: S...	8.135	28.6	ng/ul	276956	1	8.194	193936	20.0
Benzene, 1,2,3-...	8.299	28.9	ng/ul	280704	1	8.194	193936	20.0
4-Methyl-3-hept...	8.746	41.3	ng/ul	400854	1	8.194	193936	20.0
Benzene, 1,2-di...	8.822	64.0	ng/ul	620115	1	8.194	193936	20.0
Benzene, 1-ethy...	9.286	36.5	ng/ul	354034	1	8.194	193936	20.0
unknown-02	9.539	19.9	ng/ul	193133	1	8.194	193936	20.0
Benzene, 2-ethy...	9.633	18.3	ng/ul	267399	2	11.025	292642	20.0
(DEL) Alkane: S...	10.397	19.9	ng/ul	291927	2	11.025	292642	20.0
(DEL) Alkane: S...	10.949	60.4	ng/ul	884348	2	11.025	292642	20.0
(DEL) Alkane: S...	11.131	15.7	ng/ul	230143	2	11.025	292642	20.0
(DEL) Alkane: S...	11.995	17.6	ng/ul	256978	2	11.025	292642	20.0
(DEL) Alkane: S...	12.389	53.1	ng/ul	777637	2	11.025	292642	20.0
Benzenamine, 2,...	13.276	18.6	ng/ul	316992	3	14.827	340591	20.0
(DEL) Alkane: S...	13.329	11.0	ng/ul	187573	3	14.827	340591	20.0
Naphthalene, 2,...	14.004	23.9	ng/ul	407212	3	14.827	340591	20.0
Naphthalene, 2,...	14.145	28.4	ng/ul	484553	3	14.827	340591	20.0
Benzenamine, 3,...	14.175	31.1	ng/ul	529639	3	14.827	340591	20.0
3,5-Dimethyl-4-...	14.263	18.2	ng/ul	309665	3	14.827	340591	20.0
(DEL) Alkane: S...	14.680	53.1	ng/ul	904495	3	14.827	340591	20.0
Naphthalene, 1,...	15.291	20.2	ng/ul	343844	3	14.827	340591	20.0
(DEL) Alkane: S...	16.496	78.3	ng/ul	1550930	4	17.583	396346	20.0
Benzene, (1-met...	16.625	17.6	ng/ul	348821	4	17.583	396346	20.0
(DEL) Alkane: S...	17.283	35.3	ng/ul	698547	4	17.583	396346	20.0
(DEL) Alkane: S...	17.336	16.3	ng/ul	322329	4	17.583	396346	20.0
(DEL) Alkane: S...	18.023	35.1	ng/ul	695918	4	17.583	396346	20.0
n-Hexadecanoic ...	18.452	34.1	ng/ul	676355	4	17.583	396346	20.0
(DEL) Alkane: S...	18.711	22.0	ng/ul	436117	4	17.583	396346	20.0
(DEL) Alkane: S...	19.339	40.0	ng/ul	792230	4	17.583	396346	20.0
cyclohexane, [1...	19.539	17.9	ng/ul	354883	4	17.583	396346	20.0
Tridecane, 1-iodo-	19.909	17.3	ng/ul	372884	5	21.883	431232	20.0
1,2-Benzenedica...	19.939	23.8	ng/ul	512652	5	21.883	431232	20.0
(DEL) Alkane: S...	20.438	31.6	ng/ul	681023	5	21.883	431232	20.0
(DEL) Alkane: S...	20.937	27.4	ng/ul	591229	5	21.883	431232	20.0
Phosphoric acid...	21.196	32.6	ng/ul	703489	5	21.883	431232	20.0
(DEL) Alkane: S...	21.460	62.0	ng/ul	1336590	5	21.883	431232	20.0
(DEL) Alkane: S...	22.024	52.5	ng/ul	1132070	5	21.883	431232	20.0
Octane, 1,1'-ox...	22.530	100.0	ng/ul	2157010	5	21.883	431232	20.0
(DEL) Alkane: S...	22.659	52.8	ng/ul	1138360	5	21.883	431232	20.0
(DEL) Alkane: S...	23.158	20.3	ng/ul	437064	5	21.883	431232	20.0
(DEL) Alkane: S...	23.382	44.4	ng/ul	958119	5	21.883	431232	20.0
Phthalic acid, ...	23.881	20.9	ng/ul	472290	6	25.291	452308	20.0
(DEL) Alkane: S...	24.228	28.8	ng/ul	651197	6	25.291	452308	20.0
1,2-Benzenedica...	24.633	71.1	ng/ul	1608510	6	25.291	452308	20.0

1-Hexadecanethiol	24.909	46.0	ng/ul	1039500	6	25.291	452308	20.0
(DEL) Alkane: S...	25.221	20.0	ng/ul	452308	6	25.291	452308	20.0
(DEL) Alkane: S...	26.413	24.8	ng/ul	560618	6	25.291	452308	20.0
1,2-Benzenedica...	26.913	42.4	ng/ul	958311	6	25.291	452308	20.0
(DEL) Alkane: S...	27.113	53.1	ng/ul	1200350	6	25.291	452308	20.0

Instrument :
BNA_G
ClientSampleId :
BGKQ1DL